

SITE INSPECTION REPORT  
FOR  
MOBIL CHEMICAL PHOSPHORUS DIVISION  
GARY, IN  
IND000606731  
F05-8702-172

EPA Region 5 Records Ctr.



322720

PAN: FIN0110SI

JUNE 12, 1987

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**SITE INSPECTION MEMO**

**1**

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**2070 - 13 FORM**

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**SITE MAPS**

**3**

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**SITE PHOTOGRAPHS**

**4**

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**ANALYTICAL DATA**

**5**

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**1**

# 00873 NB



## ecology and environment, inc.

111 WEST JACKSON BLVD., CHICAGO, ILLINOIS 60604, TEL. 312-663-9415

International Specialists in the Environment

### M E M O R A N D U M

DATE: June 12, 1987  
TO: File  
FROM: Donald L. Clark *DLK*  
SUBJECT: Indiana/F05-8702-172/FIN0110SI  
Gary/Mobil Chemical Co. Phosphorus Division  
IND000606731

The Mobil Chemical Co. Phosphorus Division site is located in Gary, Indiana at 1040 Michigan Street, approximately 2,000 feet west of the Interstate 90-65 interchange. The property in question is now owned by Montgomery Tank Lines, which leases part of the site to Gary Products Corporation. Montgomery Tank Lines presently uses part of the facility as a transfer station for trucks carrying food products and/or soap. Gary Products is involved in soap manufacturing and packaging. Mobil operated the facility from 1965-1975, and stored phosphorus in tanks located on-site. Mobil instituted a clean-up program of tanks and tank areas before the facility was sold. A concrete catch basin, used as primary separator, is also present on-site. The site was assigned to Ecology and Environment, Inc. (FIT) for a site inspection by the U.S. EPA after being identified by Mobil Chemical Co. through EPA Form 8900-1 "Notification of Hazardous Waste Site", and a Preliminary Assessment conducted by the U.S. EPA.

Ecology and Environment, Inc. (FIT) conducted an inspection of the Mobil Chemical Co. Phosphorus Division site on February 18, 1987. During the site inspection, FIT interviewed the owner and operator and collected three on-site soil/sediment samples and a background soil sample. The FIT team expected to find a lagoon on-site, but none was present, so composite sampling around the two tank areas and within the catch basin was done. Aside from changing potential

sampling locations, the work plan was followed. Sampling was done to determine waste characteristics.

Analytical results from the soil/sediment samples indicate that hazardous substances are present in all of the site samples. Compounds and elements that were present in concentrations which can be used for HRS scoring purposes are listed below. Organic sample results for sample S2 (catch basin) show values for toluene (15,000 ppb), ethyl benzene (33,000 ppb), total xylenes (160,000 ppb), 1,2 dichlorobenzene (11,000 ppb), naphthalene (20,000 ppb), 2-methylnaphthalene (10,000 ppb), 3-nitroaniline (30,000 ppb), n-nitrosodiphenylamine (20,000 ppb), pentachlorophenol (55,000 ppb), and bis-(2-ethylhexyl)phthalate (100,000 ppb). Inorganic results for sample S2 show cadmium at 3.7 ppm and mercury at 7.0 ppm. For sample S3 (north tanks area), bis-(2-ethylhexyl)phthalate was detected at 14,000 ppb. It should also be noted that the U.S. EPA sampled insulation on pipes leading to tanks at the site to check for asbestos on 8/19/80. Sample results indicate asbestos fibers are present in the insulation. This insulation was badly weathered and peeled. FIT also found the insulation in a similar condition.

The potential for groundwater contamination exists at the site since the water table is high and surface soils are contaminated. However, no target population for groundwater exists within three miles of the site. Surface water could potentially be contaminated through surface runoff to a nearby unnamed lake. The lake is used for fishing. Contaminants may also be entering the sewer system via the catch basin. The local waste water treatment plant has detected high levels of phosphorus and has contacted Gary Products Co.. The drinking water for the area is supplied by Lake Michigan. Minimal potential for direct contact with areas of contamination exists as the site is fenced and locked at night.





POTENTIAL HAZARDOUS WASTE SITE  
SITE INSPECTION REPORT  
PART 1 - SITE LOCATION AND INSPECTION INFORMATION

I IDENTIFICATION  
01 STATE **IN** 02 SITE NUMBER **IND000606731**

II. SITE NAME AND LOCATION

01 SITE NAME (i.e., EPA common or descriptive name of site) **Mobil Chemical Co. Phosphorus Division** 02 STREET, ROUTE NO. OR SPECIFIC LOCATION IDENTIFIER **1040 Michigan St.**  
03 CITY **Gary** 04 STATE **IN** 05 ZIP CODE **46403** 06 COUNTY **Lake** 07 COUNTY CODE **089** 08 CONG. DIST. **01**

09 COORDINATES  
LATITUDE **41° 35' 30.0" N** LONGITUDE **087° 18' 45.0" W**  
10 TYPE OF OWNERSHIP (check one):  
 A PRIVATE  B FEDERAL  C STATE  D COUNTY  E MUNICIPAL  F OTHER  G UNKNOWN

III. INSPECTION INFORMATION

01 DATE OF INSPECTION **2, 18, 87** 02 SITE STATUS  ACTIVE  INACTIVE  
03 YEARS OF OPERATION **1964, 1975** — UNKNOWN  
BEGINNING YEAR ENDING YEAR

04 AGENCY PERFORMING INSPECTION (check all that apply):  
 A EPA  B EPA CONTRACTOR **Ecology & Environment, Inc.**  C MUNICIPAL  D MUNICIPAL CONTRACTOR  
 E STATE  F STATE CONTRACTOR  G OTHER

05 CHIEF INSPECTOR	06 TITLE	07 ORGANIZATION	08 TELEPHONE NO.
<b>Julie E. Kaiser</b>	<b>Biologist</b>	<b>Ecology &amp; Environment</b>	<b>(312) 663-9415</b>
09 OTHER INSPECTORS	10 TITLE	11 ORGANIZATION	12 TELEPHONE NO.
<b>Craig Almanza</b>	<b>Technician</b>	<b>Ecology &amp; Environment</b>	<b>(312) 663-9415</b>
<b>Margie Hein</b>	<b>Biologist</b>	<b>Ecology &amp; Environment</b>	<b>(312) 663-9415</b>
<b>Tom Kouris</b>	<b>Civil Engineer</b>	<b>Ecology &amp; Environment</b>	<b>(312) 663-9415</b>
			( )
			( )

13 SITE REPRESENTATIVES INTERVIEWED	14 TITLE	15 ADDRESS	16 TELEPHONE NO.
<b>Bill Keagle</b>	<b>Owner</b>	<b>Gary Products, Inc. 1040 Michigan Ave.</b>	<b>(219) 885-1644</b>
<b>Merv Bridges</b>	<b>Operations Manager</b>	<b>Montgomery Tank Lines 759 1/2 Melton Road Gary, IN 46403</b>	( )
			( )
			( )
			( )
			( )

17 ACCESS GAINED BY (check one):  PERMISSION  WARRANT  
18 TIME OF INSPECTION **10:45 a.m.** 19 WEATHER CONDITIONS **Clear, 37°F**

IV. INFORMATION AVAILABLE FROM

01 CONTACT **Dr. Don Josif** 02 OF (Agency/Organization) **U.S. EPA** 03 TELEPHONE NO. **(312) 886-0393**  
04 PERSON RESPONSIBLE FOR SITE INSPECTION FORM **Donald L. Clark** 05 AGENCY **U.S. EPA/FIT** 06 ORGANIZATION **Ecology & Environment** 07 TELEPHONE NO. **(312) 663-9415** 08 DATE **6, 15, 87**  
MONTH DAY YEAR



POTENTIAL HAZARDOUS WASTE SITE  
SITE INSPECTION REPORT  
PART 2 - WASTE INFORMATION

I IDENTIFICATION  
01 STATE: IN 02 SITE NUMBER: IN D000606731

II. WASTE STATES, QUANTITIES, AND CHARACTERISTICS

<p>01 PHYSICAL STATES (check all that apply):</p> <p><input checked="" type="checkbox"/> A SOLID <input type="checkbox"/> B POWDER FINES <input type="checkbox"/> C SLUDGE <input type="checkbox"/> D OTHER _____ (Specify)</p> <p><input type="checkbox"/> E SLURRY <input type="checkbox"/> F LIQUID <input type="checkbox"/> G GAS</p>	<p>02 WASTE QUANTITY AT SITE (Measures of waste quantities must be independent)</p> <p>TONS <u>Unknown</u> CUBIC YARDS <u>Unknown</u> NO OF DRUMS <u>1</u></p>	<p>03 WASTE CHARACTERISTICS (check all that apply):</p> <p><input type="checkbox"/> A TOXIC <input type="checkbox"/> B CORROSIVE <input type="checkbox"/> C RADIOACTIVE <input type="checkbox"/> D PERSISTENT</p> <p><input type="checkbox"/> E SOLUBLE <input type="checkbox"/> F INFECTIOUS <input type="checkbox"/> G FLAMMABLE <input type="checkbox"/> H IGNITABLE</p> <p><input type="checkbox"/> I HIGHLY VOLATILE <input type="checkbox"/> J EXPLOSIVE <input type="checkbox"/> K REACTIVE <input type="checkbox"/> L INCOMPATIBLE <input type="checkbox"/> M NOT APPLICABLE</p>
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III. WASTE TYPE

CATEGORY	SUBSTANCE NAME	01 GROSS AMOUNT	02 UNIT OF MEASURE	03 COMMENTS
	Misc.	unknown		P4, Asbestos
OLW	OILY WASTE			
SOL	SOLVENTS			
PSD	PESTICIDES			
OCC	OTHER ORGANIC CHEMICALS			
IOC	INORGANIC CHEMICALS	unknown		salts
ACD	ACIDS	unknown		pH < 3
BAS	BASES			
MES	HEAVY METALS	unknown		arsenic, antimony, selenium, iron, manganese, magnesium

IV. HAZARDOUS SUBSTANCES (See Appendix for most frequently cited CAS Numbers)

01 CATEGORY	02 SUBSTANCE NAME	03 CAS NUMBER	04 STORAGE/ DISPOSAL METHOD	05 CONCENTRATION	06 MEASURE OF CONCENTRATION
The following hazardous substances were detected in soil/sediment samples collected at the site, and can be used for HRS scoring:					
Organics	EJ 105 toluene			15,000	ug/Kg
	(Catch Basin) ethyl benzene			33,000	
	total xylenes			160,000	
	1,2 dichlorobenzene			11,000	
	naphthalene			20,000	
	2-methyl-naphthalene			10,000	
	3-nitroaniline			30,000	
	n-nitrosodiphenylamine			20,000	
	pentachlorophenol			55,000	
	bis-(2-ethyl hexyl) phthalate			100,000	
	EJ 106 bis-(2-ethyl hexyl) phthalate			14,000	
	(North Tanks)				
Inorganics	MEI 376 Cadmium			3.7	mg/Kg
	(Catch Basin) Mercury			7.0	

V. FEEDSTOCKS (See Appendix for CAS Numbers)

CATEGORY	01 FEEDSTOCK NAME	02 CAS NUMBER	CATEGORY	01 FEEDSTOCK NAME	02 CAS NUMBER
FDS	None		FDS		
FDS			FDS		
FDS			FDS		
FDS			FDS		

VI. SOURCES OF INFORMATION (See specific references e.g. state files, sample analysis reports)

Attachment to EPA Form 8900-1 "Notification of Hazardous Waste Site" - Form B, and letter from W.R. Smithy, Jr. (U.S. EPA) June 9, 1981  
Sample analysis reports for Mobil Chem. Phosph. Div., E4E/FIT site inspection.



**POTENTIAL HAZARDOUS WASTE SITE  
SITE INSPECTION REPORT**  
PART 3 - DESCRIPTION OF HAZARDOUS CONDITIONS AND INCIDENTS

I. IDENTIFICATION	
01 STATE <b>IN</b>	02 SITE NUMBER <b>IND000606731</b>

**II. HAZARDOUS CONDITIONS AND INCIDENTS**

01  A. GROUNDWATER CONTAMINATION      02  OBSERVED (DATE: \_\_\_\_\_)       POTENTIAL       ALLEGED  
03 POPULATION POTENTIALLY AFFECTED: -0-      04 NARRATIVE DESCRIPTION  
High potential for groundwater contamination exists since the water table is found within 15 feet of the ground surface and surface soils are contaminated. No target population for groundwater exists.

01  B. SURFACE WATER CONTAMINATION      02  OBSERVED (DATE: \_\_\_\_\_)       POTENTIAL       ALLEGED  
03 POPULATION POTENTIALLY AFFECTED: \_\_\_\_\_      04 NARRATIVE DESCRIPTION  
Potential for surface water contamination exists through the possibility of contaminants entering nearby lake via surface runoff. Lake is unnamed but on-site representatives have seen people fishing in it.

01  C. CONTAMINATION OF AIR      02  OBSERVED (DATE: \_\_\_\_\_)       POTENTIAL       ALLEGED  
03 POPULATION POTENTIALLY AFFECTED: \_\_\_\_\_      04 NARRATIVE DESCRIPTION  
Insulation found on pipes leading to tanks is exposed and damaged, and is asbestos containing, therefore a potential hazard exists.

01  D. FIRE/EXPLOSIVE CONDITIONS      02  OBSERVED (DATE: \_\_\_\_\_)       POTENTIAL       ALLEGED  
03 POPULATION POTENTIALLY AFFECTED: 12      04 NARRATIVE DESCRIPTION  
A site representative reports that phosphorus ignited on contact with air when he was taking valves apart during a clean-up process.

01  E. DIRECT CONTACT      02  OBSERVED (DATE: \_\_\_\_\_)       POTENTIAL       ALLEGED  
03 POPULATION POTENTIALLY AFFECTED: -0-      04 NARRATIVE DESCRIPTION  
The site is surrounded by fencing and the front gate is locked at night; therefore, the potential for an incident via direct contact is low. The catch basin is not covered.

01  F. CONTAMINATION OF SOIL      02  OBSERVED (DATE: \_\_\_\_\_)       POTENTIAL       ALLEGED  
03 AREA POTENTIALLY AFFECTED: 5.3      04 NARRATIVE DESCRIPTION  
(Acres)  
Composite soil samples taken near the north and south tanks show contamination.

01  G. DRINKING WATER CONTAMINATION      02  OBSERVED (DATE: \_\_\_\_\_)       POTENTIAL       ALLEGED  
03 POPULATION POTENTIALLY AFFECTED: -0-      04 NARRATIVE DESCRIPTION  
N/A - Drinking water source for all populations within a 3-mile radius is Lake Michigan to which there is no surface water route attributable to the site.

01  H. WORKER EXPOSURE/INJURY      02  OBSERVED (DATE: \_\_\_\_\_)       POTENTIAL       ALLEGED  
03 WORKERS POTENTIALLY AFFECTED: 12      04 NARRATIVE DESCRIPTION  
Owner was cutting beams with a welder and felt ill afterwards. Test results were inconclusive. Possibly hydrogen gas was produced from a reaction with phosphorus.

01  I. POPULATION EXPOSURE/INJURY      02  OBSERVED (DATE: \_\_\_\_\_)       POTENTIAL       ALLEGED  
03 POPULATION POTENTIALLY AFFECTED: -0-      04 NARRATIVE DESCRIPTION  
There is minimal potential for population exposure/injury as site area and catch basin are fenced with good security.



POTENTIAL HAZARDOUS WASTE SITE  
SITE INSPECTION REPORT  
PART 3 - DESCRIPTION OF HAZARDOUS CONDITIONS AND INCIDENTS

I. IDENTIFICATION  
01 STATE IN 02 SITE NUMBER IND000606731

II. HAZARDOUS CONDITIONS AND INCIDENTS (Continued)

01  J. DAMAGE TO FLORA 02  OBSERVED (DATE: \_\_\_\_\_)  POTENTIAL  ALLEGED  
04 NARRATIVE DESCRIPTION

There have been no signs of damage to flora.

01  K. DAMAGE TO FAUNA 02  OBSERVED (DATE: \_\_\_\_\_)  POTENTIAL  ALLEGED  
04 NARRATIVE DESCRIPTION (include name(s) of species)

There have been no signs of damage to fauna.

01  L. CONTAMINATION OF FOOD CHAIN 02  OBSERVED (DATE: \_\_\_\_\_)  POTENTIAL  ALLEGED  
04 NARRATIVE DESCRIPTION

See J and K above.

01  M. UNSTABLE CONTAINMENT OF WASTES 02  OBSERVED (DATE: \_\_\_\_\_)  POTENTIAL  ALLEGED  
(Spills/Runoff/Standing liquids, Leaking drums)

03 POPULATION POTENTIALLY AFFECTED: 12+ 04 NARRATIVE DESCRIPTION  
Contaminants in soils indicate possible tank leakage or spills. Water in catch basin flows into sewer system.

01  N. DAMAGE TO OFFSITE PROPERTY 02  OBSERVED (DATE: \_\_\_\_\_)  POTENTIAL  ALLEGED  
04 NARRATIVE DESCRIPTION

See B, J, and K above.

01  O. CONTAMINATION OF SEWERS, STORM DRAINS, WWTPs 02  OBSERVED (DATE: \_\_\_\_\_)  POTENTIAL  ALLEGED  
04 NARRATIVE DESCRIPTION

The catch basin leads to the city's sewer system. WWTP has picked up phosphorus in their analyses on several occasions and has notified the facility. Other contaminants found in catch basin sediments could be entering sewer system.

01  P. ILLEGAL/UNAUTHORIZED DUMPING 02  OBSERVED (DATE: \_\_\_\_\_)  POTENTIAL  ALLEGED  
04 NARRATIVE DESCRIPTION

There have been no incidents on site.

05 DESCRIPTION OF ANY OTHER KNOWN, POTENTIAL, OR ALLEGED HAZARDS

None

III. TOTAL POPULATION POTENTIALLY AFFECTED: 12

IV. COMMENTS

Potential for contamination of air cannot currently be substantiated, since no air samples have been collected.

V. SOURCES OF INFORMATION (Cite specific references, e.g., state files, sample analysis, reports)

FIT Site Inspection 2/12/87



POTENTIAL HAZARDOUS WASTE SITE  
SITE INSPECTION  
PART 4 - PERMIT AND DESCRIPTIVE INFORMATION

I. IDENTIFICATION  
01 STATE 02 SITE NUMBER  
IN IND000606731

II. PERMIT INFORMATION

01 TYPE OF PERMIT ISSUED (Check all that apply)	02 PERMIT NUMBER	03 DATE ISSUED	04 EXPIRATION DATE	05 COMMENTS
<input type="checkbox"/> A. NPDES				
<input type="checkbox"/> B. UIC				
<input type="checkbox"/> C. AIR				
<input type="checkbox"/> D. RCRA				
<input type="checkbox"/> E. RCRA INTERIM STATUS				
<input type="checkbox"/> F. SPCC PLAN				
<input type="checkbox"/> G. STATE (Specify)				
<input type="checkbox"/> H. LOCAL (Specify)				
<input type="checkbox"/> I. OTHER (Specify)				
<input checked="" type="checkbox"/> J. NONE				

III. SITE DESCRIPTION

01 STORAGE/DISPOSAL (Check all that apply)	02 AMOUNT	03 UNIT OF MEASURE	04 TREATMENT (Check all that apply)	05 OTHER
<input type="checkbox"/> A. SURFACE IMPOUNDMENT <input type="checkbox"/> B. PILES <input type="checkbox"/> C. DRUMS, ABOVE GROUND <input type="checkbox"/> D. TANK, ABOVE GROUND <input type="checkbox"/> E. TANK, BELOW GROUND <input type="checkbox"/> F. LANDFILL <input type="checkbox"/> G. LANDFARM <input type="checkbox"/> H. OPEN DUMP <input checked="" type="checkbox"/> I. OTHER <u>Catch Basin</u> <small>(Specify)</small>			<input type="checkbox"/> A. INCENERATION <input type="checkbox"/> B. UNDERGROUND INJECTION <input type="checkbox"/> C. CHEMICAL/PHYSICAL <input type="checkbox"/> D. BIOLOGICAL <input type="checkbox"/> E. WASTE OIL PROCESSING <input type="checkbox"/> F. SOLVENT RECOVERY <input type="checkbox"/> G. OTHER RECYCLING/RECOVERY <input checked="" type="checkbox"/> H. OTHER <u>Neutralization</u> <small>(Specify)</small>	<input checked="" type="checkbox"/> A. BUILDINGS ON SITE  4  06 AREA OF SITE  5.3 (Acres)

07 COMMENTS

No permits are known to exist.

IV. CONTAINMENT

01 CONTAINMENT OF WASTES (Check one)

A. ADEQUATE, SECURE       B. MODERATE       C. INADEQUATE, POOR       D. INSECURE, UNSOUND, DANGEROUS

02 DESCRIPTION OF DRUMS, DIKING, LINERS, BARRIERS, ETC.  
 Storage tanks appear in sound condition. However, soil contamination indicates containment problems. Bottom and walls of catch basin are concrete, but effluent flows into sewer system through piping. Uncertain as to whether a landfill exists on site. There may have been some dumping into an on-site landfill prior to 1975.

V. ACCESSIBILITY

01 WASTE EASILY ACCESSIBLE:  YES  NO

02 COMMENTS

The entire site is surrounded by a fence, and the front gate is locked at night.

VI. SOURCES OF INFORMATION (Cite specific references, e.g. state files, sample analysis, reports)

Attachment to EPA Form 8900-1 "Notification of Hazardous Waste Site", 6/1/81  
 FIT Site Inspection 2/18/87



POTENTIAL HAZARDOUS WASTE SITE  
SITE INSPECTION REPORT  
PART 5 - WATER, DEMOGRAPHIC, AND ENVIRONMENTAL DATA

I. IDENTIFICATION  
01 STATE **IN** 02 SITE NUMBER **IND 00060673**

II. DRINKING WATER SUPPLY

01 TYPE OF DRINKING SUPPLY <small>(Check as applicable)</small>	SURFACE		WELL		02 STATUS			03 DISTANCE TO SITE	
	COMMUNITY	A <input checked="" type="checkbox"/>	B <input type="checkbox"/>	ENDANGERED	AFFECTED	MONITORED	A	<u>&gt;3</u>	(mi)
NON-COMMUNITY	C <input checked="" type="checkbox"/>	D <input type="checkbox"/>	D <input type="checkbox"/>	E <input type="checkbox"/>	F <input checked="" type="checkbox"/>	B	<u>&gt;3</u>	(mi)	

III. GROUNDWATER

01 GROUNDWATER USE IN VICINITY (Check one)

A ONLY SOURCE FOR DRINKING     B DRINKING (Other sources available)  
COMMERCIAL INDUSTRIAL IRRIGATION (No other water sources available)

C COMMERCIAL INDUSTRIAL IRRIGATION (Limit other sources available)     D NOT USED, UNUSABLE

02 POPULATION SERVED BY GROUND WATER -0-

03 DISTANCE TO NEAREST DRINKING WATER WELL >3 (mi)

04 DEPTH TO GROUNDWATER <15 (ft)

05 DIRECTION OF GROUNDWATER FLOW NNE

06 DEPTH TO AQUIFER OF CONCERN <15 (ft)

07 POTENTIAL YIELD OF AQUIFER Unknown (gpd)

08 SOLE SOURCE AQUIFER  YES  NO

09 DESCRIPTION OF WELLS (including usage, depth, and location relative to population and buildings)  
*There are no groundwater wells known to exist within a three-mile radius of the site. All water is supplied by Lake Michigan.*

10 RECHARGE AREA	COMMENTS	11 DISCHARGE AREA	COMMENTS
<input checked="" type="checkbox"/> YES <input type="checkbox"/> NO	<i>site lies in the possible recharge area of the Grand and Little Calumet Rivers.</i>	<input checked="" type="checkbox"/> YES <input type="checkbox"/> NO	<i>Possibly a local discharge area in the vicinity of the small unnamed lake west of the site.</i>

IV. SURFACE WATER

01 SURFACE WATER USE (Check one)

A. RESERVOIR, RECREATION DRINKING WATER SOURCE     B. IRRIGATION, ECONOMICALLY IMPORTANT RESOURCES     C. COMMERCIAL, INDUSTRIAL     D. NOT CURRENTLY USED

02 AFFECTED/POTENTIALLY AFFECTED BODIES OF WATER

NAME	AFFECTED	DISTANCE TO SITE
<i>Small unnamed lake located west of facility</i>	<input type="checkbox"/>	<u>500</u> ft.
<i>Grand Calumet River</i>	<input type="checkbox"/>	<u>5540</u> (mi)
<i>Little Calumet River</i>	<input type="checkbox"/>	<u>7290</u> (mi)

V. DEMOGRAPHIC AND PROPERTY INFORMATION

01 TOTAL POPULATION WITHIN

ONE (1) MILE OF SITE    TWO (2) MILES OF SITE    THREE (3) MILES OF SITE

A. ~20,000 NO. OF PERSONS    B. ~55,000 NO. OF PERSONS    C. ~114,000 NO. OF PERSONS

02 DISTANCE TO NEAREST POPULATION 1690 ft.

03 NUMBER OF BUILDINGS WITHIN TWO (2) MILES OF SITE 14,800

04 DISTANCE TO NEAREST OFF-SITE BUILDING 300 ft.

05 POPULATION WITHIN VICINITY OF SITE (Provide narrative description of nature of population within vicinity of site e.g. rural village densely populated urban area)  
*Densely populated urban area exists around the site, as well as highly industrialized areas nearby.*



POTENTIAL HAZARDOUS WASTE SITE  
SITE INSPECTION REPORT  
PART 5 - WATER, DEMOGRAPHIC, AND ENVIRONMENTAL DATA

I. IDENTIFICATION  
01 STATE **IN** 02 SITE NUMBER **IND000606731**

VI. ENVIRONMENTAL INFORMATION

01 PERMEABILITY OF UNSATURATED ZONE (Check one)

A  $10^{-8} - 10^{-9}$  cm/sec  B  $10^{-6} - 10^{-7}$  cm/sec  C  $10^{-4} - 10^{-5}$  cm/sec  D GREATER THAN  $10^{-3}$  cm/sec

02 PERMEABILITY OF BEDROCK (Check one)

A IMPERMEABLE (Less than  $10^{-6}$  cm/sec)  B RELATIVELY IMPERMEABLE ( $10^{-4} - 10^{-6}$  cm/sec)  C RELATIVELY PERMEABLE ( $10^{-2} - 10^{-4}$  cm/sec)  D VERY PERMEABLE (Greater than  $10^{-2}$  cm/sec)

03 DEPTH TO BEDROCK

~ 160 (ft)

04 DEPTH OF CONTAMINATED SOIL ZONE

N/A (ft)

05 SOIL pH

7.75

06 NET PRECIPITATION

6 (in)

07 ONE YEAR 24 HOUR RAINFALL

2.4 (in)

08 SLOPE

SITE SLOPE 43 %

DIRECTION OF SITE SLOPE

N/A

TERRAIN AVERAGE SLOPE

43 %

09 FLOOD POTENTIAL

SITE IS IN N/A YEAR FLOODPLAIN

10

N/A

SITE IS ON BARRIER ISLAND, COASTAL HIGH HAZARD AREA, RIVERINE FLOODWAY

11 DISTANCE TO WETLANDS (5 acre minimum)

ESTUARINE

A N/A (mi)

OTHER

B 4650 (mit ft)

12 DISTANCE TO CRITICAL HABITAT (of endangered species)

> 3 (mi)

ENDANGERED SPECIES N/A

13 LAND USE IN VICINITY

DISTANCE TO

COMMERCIAL/INDUSTRIAL

A 1200 (mit ft)

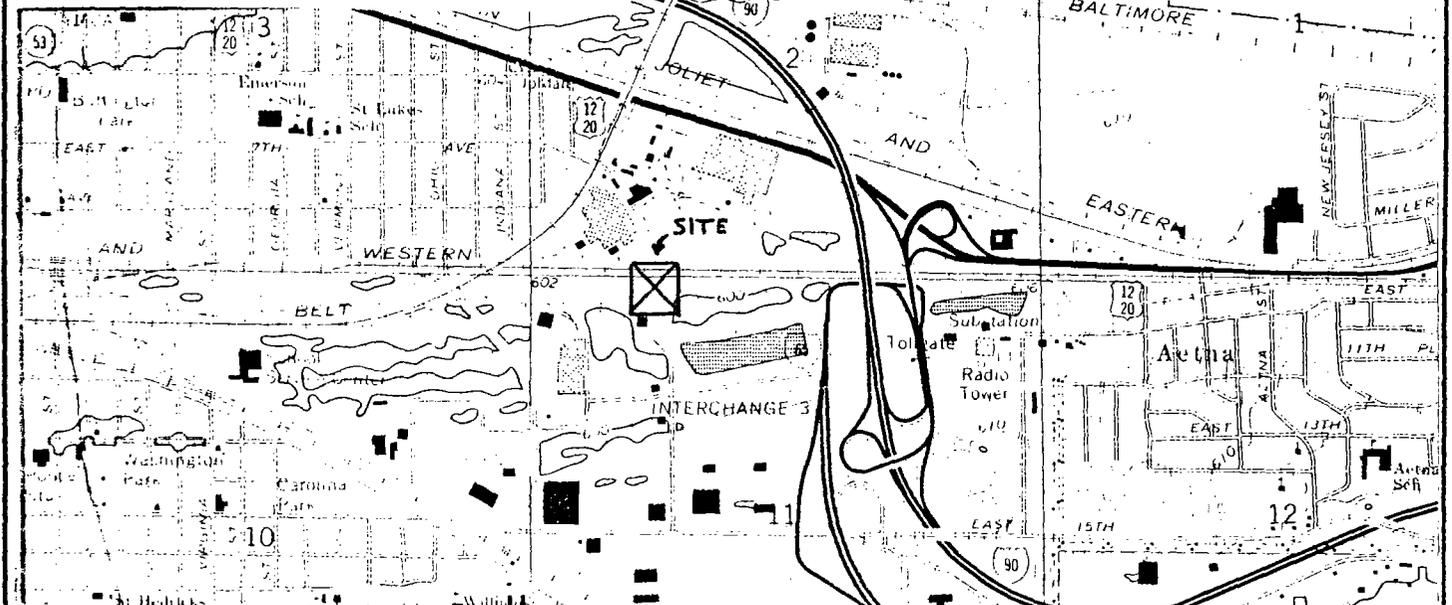
RESIDENTIAL AREAS, NATIONAL/STATE PARKS, FORESTS, OR WILDLIFE RESERVES

B 1850 (mit ft)

AGRICULTURAL LANDS  
PRIME AG LAND AG LAND

C > 3 (mi) D > 3 (mi)

14 DESCRIPTION OF SITE IN RELATION TO SURROUNDING TOPOGRAPHY



VII. SOURCES OF INFORMATION (Cite specific references e.g. state fees sample analysis reports)

Environmental Geology of Lake and Porter Counties, IN - An Aid to Planning:  
IDNR Envir. Study 8.  
Freeze, R.A. and J.A. Cherry, Groundwater, Prentice Hall, NY, 1979.  
Gary, IN 7 1/2 Quad. Topographic Map (USGS)



POTENTIAL HAZARDOUS WASTE SITE  
SITE INSPECTION REPORT  
PART 6 - SAMPLE AND FIELD INFORMATION

I. IDENTIFICATION

D1 STATE: IN D2 SITE NUMBER: IND 000606731

II. SAMPLES TAKEN

SAMPLE TYPE	D1 NUMBER OF SAMPLES TAKEN	D2 SAMPLES SENT TO	D3 ESTIMATED DATE RESULTS AVAILABLE
GROUNDWATER			
SURFACE WATER			
WASTE			
AIR		Inorganics: Century Labs, Inc.	March 26, 1987
RUNOFF		1501 Grandview Ave., Thorofare, NJ 08086	
SPILL		Contact: Al Tordini	
SOIL/SEDIMENT	4	Organics: Analytical Technologies, 5550 Morehouse Dr. San Diego, CA 92121	March 30, 1987
VEGETATION			
OTHER		Contact: Bob Woods	

III. FIELD MEASUREMENTS TAKEN

D1 TYPE	D2 COMMENTS
None	

IV. PHOTOGRAPHS AND MAPS

D1 TYPE <input checked="" type="checkbox"/> GROUND <input type="checkbox"/> AERIAL	D2 IN CUSTODY OF Ecology & Environment, Inc. <small>(Name of organization or individual)</small>
D3 MAPS <input type="checkbox"/> YES <input checked="" type="checkbox"/> NO	D4 LOCATION OF MAPS N/A

V. OTHER FIELD DATA COLLECTED (Provide narrative description)

None

VI. SOURCES OF INFORMATION (Cite specific references e.g. State files sample analysis reports)

2/18/87 ETE/FIT site inspection



POTENTIAL HAZARDOUS WASTE SITE  
SITE INSPECTION REPORT  
PART 7 - OWNER INFORMATION

I IDENTIFICATION  
01 STATE 02 SITE NUMBER  
IN IND000606731

II. CURRENT OWNER(S)				PARENT COMPANY (IF APPLICABLE)			
01 NAME MONTGOMERY TANK LINES		02 D+B NUMBER unknown		08 NAME MONTGOMERY TANK LINES		09 D+B NUMBER unknown	
03 STREET ADDRESS (P.O. Box, RFD #, etc.) 1040 MICHIGAN			04 SIC CODE unknown	10 STREET ADDRESS (P.O. Box, RFD #, etc.) 3106 CENTRAL AVE. PO BOX 2628			11 SIC CODE unknown
05 CITY GARY		06 STATE IN	07 ZIP CODE 46403	12 CITY PLANT CITY		13 STATE FL	14 ZIP CODE 33566
01 NAME		02 D+B NUMBER		08 NAME		09 D+B NUMBER	
03 STREET ADDRESS (P.O. Box, RFD #, etc.)			04 SIC CODE	10 STREET ADDRESS (P.O. Box, RFD #, etc.)			11 SIC CODE
05 CITY		06 STATE	07 ZIP CODE	12 CITY		13 STATE	14 ZIP CODE
01 NAME		02 D+B NUMBER		08 NAME		09 D+B NUMBER	
03 STREET ADDRESS (P.O. Box, RFD #, etc.)			04 SIC CODE	10 STREET ADDRESS (P.O. Box, RFD #, etc.)			11 SIC CODE
05 CITY		06 STATE	07 ZIP CODE	12 CITY		13 STATE	14 ZIP CODE
01 NAME		02 D+B NUMBER		08 NAME		09 D+B NUMBER	
03 STREET ADDRESS (P.O. Box, RFD #, etc.)			04 SIC CODE	10 STREET ADDRESS (P.O. Box, RFD #, etc.)			11 SIC CODE
05 CITY		06 STATE	07 ZIP CODE	12 CITY		13 STATE	14 ZIP CODE

III. PREVIOUS OWNER(S) (LIST MOST RECENT FIRST)				IV. PARENT COMPANY OF PREVIOUS OWNER			
01 NAME MOBIL CHEMICAL - PHOSPHORUS DIVISION		02 D+B NUMBER unknown		01 NAME MOBIL OIL CORP.		02 D+B NUMBER unknown	
03 STREET ADDRESS (P.O. Box, RFD #, etc.) P.O. Box 26683			04 SIC CODE unknown	03 STREET ADDRESS (P.O. Box, RFD #, etc.) 150 E. 42nd St.			04 SIC CODE unknown
05 CITY RICHMOND		06 STATE VA	07 ZIP CODE 23261	05 CITY NEW YORK		06 STATE NY	07 ZIP CODE 10017
01 NAME		02 D+B NUMBER		01 NAME		02 D+B NUMBER	
03 STREET ADDRESS (P.O. Box, RFD #, etc.)			04 SIC CODE	03 STREET ADDRESS (P.O. Box, RFD #, etc.)			04 SIC CODE
05 CITY		06 STATE	07 ZIP CODE	05 CITY		06 STATE	07 ZIP CODE
01 NAME		02 D+B NUMBER		01 NAME		02 D+B NUMBER	
03 STREET ADDRESS (P.O. Box, RFD #, etc.)			04 SIC CODE	03 STREET ADDRESS (P.O. Box, RFD #, etc.)			04 SIC CODE
05 CITY		06 STATE	07 ZIP CODE	05 CITY		06 STATE	07 ZIP CODE

V. SOURCES OF INFORMATION (Cite specific references e.g. state files, sample analysis reports)  
E + E FIT SITE INSPECTION 2/18/87  
Attachment to EPA form 8900-1 "Notification of Hazardous Waste Site" dated 6/9/81.



POTENTIAL HAZARDOUS WASTE SITE  
SITE INSPECTION REPORT  
PART 8 - OPERATOR INFORMATION

I. IDENTIFICATION  
01 STATE 02 SITE NUMBER  
IN IND000606731

II. CURRENT OPERATOR (Provide if different from owner)				OPERATOR'S PARENT COMPANY (if applicable)			
01 NAME Gary Products Corp.		02 D+B NUMBER unknown		10 NAME		11 D+B NUMBER	
03 STREET ADDRESS (P.O. Box, RFD #, etc.) 1040 Michigan St.		04 SIC CODE unknown		12 STREET ADDRESS (P.O. Box, RFD #, etc.)		13 SIC CODE	
05 CITY Gary		06 STATE IN	07 ZIP CODE 46403	14 CITY		15 STATE	16 ZIP CODE
08 YEARS OF OPERATION	09 NAME OF OWNER						
III. PREVIOUS OPERATOR(S) (List most recent first; provide only if different from owner)				PREVIOUS OPERATORS' PARENT COMPANIES (if applicable)			
01 NAME		02 D+B NUMBER		10 NAME		11 D+B NUMBER	
03 STREET ADDRESS (P.O. Box, RFD #, etc.)		04 SIC CODE		12 STREET ADDRESS (P.O. Box, RFD #, etc.)		13 SIC CODE	
05 CITY		06 STATE	07 ZIP CODE	14 CITY		15 STATE	16 ZIP CODE
08 YEARS OF OPERATION	09 NAME OF OWNER DURING THIS PERIOD						
01 NAME		02 D+B NUMBER		10 NAME		11 D+B NUMBER	
03 STREET ADDRESS (P.O. Box, RFD #, etc.)		04 SIC CODE		12 STREET ADDRESS (P.O. Box, RFD #, etc.)		13 SIC CODE	
05 CITY		06 STATE	07 ZIP CODE	14 CITY		15 STATE	16 ZIP CODE
08 YEARS OF OPERATION	09 NAME OF OWNER DURING THIS PERIOD						
01 NAME		02 D+B NUMBER		10 NAME		11 D+B NUMBER	
03 STREET ADDRESS (P.O. Box, RFD #, etc.)		04 SIC CODE		12 STREET ADDRESS (P.O. Box, RFD #, etc.)		13 SIC CODE	
05 CITY		06 STATE	07 ZIP CODE	14 CITY		15 STATE	16 ZIP CODE
08 YEARS OF OPERATION	09 NAME OF OWNER DURING THIS PERIOD						

IV. SOURCES OF INFORMATION (Cite specific references, e.g., state files, sample analysis, reports)

FIT Site Inspection 2/18/87



POTENTIAL HAZARDOUS WASTE SITE  
SITE INSPECTION REPORT  
PART 9 - GENERATOR/TRANSPORTER INFORMATION

I. IDENTIFICATION  
01 STATE 02 SITE NUMBER  
IN IND000606731

II. ON-SITE GENERATOR (PREVIOUS)

01 NAME MOBIL CHEMICAL CO.		02 D+B NUMBER unk	
03 STREET ADDRESS (P.O. Box, RFD#, etc.) 1040 Michigan		04 SIC CODE unk	
05 CITY GARY	06 STATE IN	07 ZIP CODE 46403	

III. OFF-SITE GENERATOR(S) (PREVIOUS)

01 NAME MOBIL CHEMICAL CO.		02 D+B NUMBER unk		01 NAME		02 D+B NUMBER	
03 STREET ADDRESS (P.O. Box, RFD#, etc.) 1040 Michigan		04 SIC CODE unk		03 STREET ADDRESS (P.O. Box, RFD#, etc.)		04 SIC CODE	
05 CITY GARY	06 STATE IN	07 ZIP CODE 46403		05 CITY	06 STATE	07 ZIP CODE	
01 NAME		02 D+B NUMBER		01 NAME		02 D+B NUMBER	
03 STREET ADDRESS (P.O. Box, RFD#, etc.)		04 SIC CODE		03 STREET ADDRESS (P.O. Box, RFD#, etc.)		04 SIC CODE	
05 CITY	06 STATE	07 ZIP CODE		05 CITY	06 STATE	07 ZIP CODE	

IV. TRANSPORTER(S)

01 NAME SANITATION SERVICE		02 D+B NUMBER N/A		01 NAME CALUMET CITY		02 D+B NUMBER N/A	
03 STREET ADDRESS (P.O. Box, RFD#, etc.) UNKOWN		04 SIC CODE N/A		03 STREET ADDRESS (P.O. Box, RFD#, etc.) UNK		04 SIC CODE N/A	
05 CITY UNK	06 STATE UNK	07 ZIP CODE UNK		05 CITY GARY	06 STATE IN	07 ZIP CODE UNK	
01 NAME MOBIL CHEMICAL COMPANY		02 D+B NUMBER N/A		01 NAME		02 D+B NUMBER	
03 STREET ADDRESS (P.O. Box, RFD#, etc.) P.O. Box 26683		04 SIC CODE N/A		03 STREET ADDRESS (P.O. Box, RFD#, etc.)		04 SIC CODE	
05 CITY RICHMOND	06 STATE VA	07 ZIP CODE 23261		05 CITY	06 STATE	07 ZIP CODE	

V. SOURCES OF INFORMATION (See specific references, e.g., state files, sample analysis, reports)

ATTACHMENT TO EPA FORM 8700-1 "NOTIFICATION OF HAZARDOUS WASTE SITE"  
FORM C - HAZARDOUS WASTE



POTENTIAL HAZARDOUS WASTE SITE  
SITE INSPECTION REPORT  
PART 10 - PAST RESPONSE ACTIVITIES

I. IDENTIFICATION  
01 STATE 02 SITE NUMBER  
IN INDO00606731

II. PAST RESPONSE ACTIVITIES

01 <input type="checkbox"/> A. WATER SUPPLY CLOSED 04 DESCRIPTION N/A	02 DATE _____	03 AGENCY _____
01 <input type="checkbox"/> B. TEMPORARY WATER SUPPLY PROVIDED 04 DESCRIPTION N/A	02 DATE _____	03 AGENCY _____
01 <input type="checkbox"/> C. PERMANENT WATER SUPPLY PROVIDED 04 DESCRIPTION N/A	02 DATE _____	03 AGENCY _____
01 <input type="checkbox"/> D. SPILLED MATERIAL REMOVED 04 DESCRIPTION N/A	02 DATE _____	03 AGENCY _____
01 <input type="checkbox"/> E. CONTAMINATED SOIL REMOVED 04 DESCRIPTION N/A	02 DATE _____	03 AGENCY _____
01 <input type="checkbox"/> F. WASTE REPACKAGED 04 DESCRIPTION N/A	02 DATE _____	03 AGENCY _____
01 <input checked="" type="checkbox"/> G. WASTE DISPOSED ELSEWHERE 04 DESCRIPTION 100 Tons of waste to Gary City Landfill - mobil 1,000 gallons of meta-phosphorus to unknown company that uses phosphorus - Gary Products Corp.	02 DATE 1975	03 AGENCY _____
01 <input checked="" type="checkbox"/> H. ON SITE BURIAL 04 DESCRIPTION 100 Tons of mono-industrial waste in landfill	02 DATE unknown	03 AGENCY _____
01 <input checked="" type="checkbox"/> I. IN SITU CHEMICAL TREATMENT 04 DESCRIPTION Neutralizing treatment at site.	02 DATE unknown	03 AGENCY Mobil Chem.
01 <input type="checkbox"/> J. IN SITU BIOLOGICAL TREATMENT 04 DESCRIPTION N/A	02 DATE _____	03 AGENCY _____
01 <input type="checkbox"/> K. IN SITU PHYSICAL TREATMENT 04 DESCRIPTION N/A	02 DATE _____	03 AGENCY _____
01 <input type="checkbox"/> L. ENCAPSULATION 04 DESCRIPTION N/A	02 DATE _____	03 AGENCY _____
01 <input type="checkbox"/> M. EMERGENCY WASTE TREATMENT 04 DESCRIPTION N/A	02 DATE _____	03 AGENCY _____
01 <input type="checkbox"/> N. CUTOFF WALLS 04 DESCRIPTION N/A	02 DATE _____	03 AGENCY _____
01 <input type="checkbox"/> O. EMERGENCY DIKING/SURFACE WATER DIVERSION 04 DESCRIPTION N/A	02 DATE _____	03 AGENCY _____
01 <input type="checkbox"/> P. CUTOFF TRENCHES/SUMP 04 DESCRIPTION N/A	02 DATE _____	03 AGENCY _____
01 <input type="checkbox"/> Q. SUBSURFACE CUTOFF WALL 04 DESCRIPTION N/A	02 DATE _____	03 AGENCY _____



POTENTIAL HAZARDOUS WASTE SITE  
SITE INSPECTION REPORT  
PART 10 - PAST RESPONSE ACTIVITIES

I. IDENTIFICATION  
01 STATE | 02 SITE NUMBER  
IN | IND000606731

II PAST RESPONSE ACTIVITIES (Continued)

01 <input type="checkbox"/> R. BARRIER WALLS CONSTRUCTED 04 DESCRIPTION N/A	02 DATE _____	03 AGENCY _____
01 <input type="checkbox"/> S. CAPPING/COVERING 04 DESCRIPTION N/A	02 DATE _____	03 AGENCY _____
01 <input type="checkbox"/> T. BULK TANKAGE REPAIRED 04 DESCRIPTION Unknown	02 DATE _____	03 AGENCY _____
01 <input type="checkbox"/> U. GROUT CURTAIN CONSTRUCTED 04 DESCRIPTION N/A	02 DATE _____	03 AGENCY _____
01 <input type="checkbox"/> V. BOTTOM SEALED 04 DESCRIPTION N/A	02 DATE _____	03 AGENCY _____
01 <input type="checkbox"/> W. GAS CONTROL 04 DESCRIPTION N/A	02 DATE _____	03 AGENCY _____
01 <input type="checkbox"/> X. FIRE CONTROL 04 DESCRIPTION N/A	02 DATE _____	03 AGENCY _____
01 <input type="checkbox"/> Y. LEACHATE TREATMENT 04 DESCRIPTION N/A	02 DATE _____	03 AGENCY _____
01 <input type="checkbox"/> Z. AREA EVACUATED 04 DESCRIPTION N/A	02 DATE _____	03 AGENCY _____
01 <input type="checkbox"/> 1. ACCESS TO SITE RESTRICTED 04 DESCRIPTION N/A	02 DATE _____	03 AGENCY _____
01 <input type="checkbox"/> 2. POPULATION RELOCATED 04 DESCRIPTION N/A	02 DATE _____	03 AGENCY _____
01 <input checked="" type="checkbox"/> 3. OTHER REMEDIAL ACTIVITIES 04 DESCRIPTION Mobil's records indicate that all tanks and adjoining lines were cleaned to bare metal. Gary Products contends that the reactor was 1/2 full of meta-phosphorus and a "heel" of phosphorus was left in the tanks. Also see Sections D & H for other indications of site not being properly cleaned up by Mobil.	02 DATE _____	03 AGENCY _____

III. SOURCES OF INFORMATION (Cite specific references, e.g., state files, sample analysis, reports)

FIT Site Inspection 2/18/87  
Attachment to EPA Form 8900-1 "Notification of Hazardous Waste Site"



POTENTIAL HAZARDOUS WASTE SITE  
SITE INSPECTION REPORT  
PART 11 - ENFORCEMENT INFORMATION

I. IDENTIFICATION	
01 STATE	02 SITE NUMBER
IN	IND00606731

II. ENFORCEMENT INFORMATION

01 PAST REGULATORY/ENFORCEMENT ACTION  YES  NO

02 DESCRIPTION OF FEDERAL, STATE, LOCAL REGULATORY/ENFORCEMENT ACTION

*There is no indication that any enforcement action was taken against Mobil Chemicals at this site.*

III. SOURCES OF INFORMATION (Cite specific references, e.g., state files, sample analysis, reports)

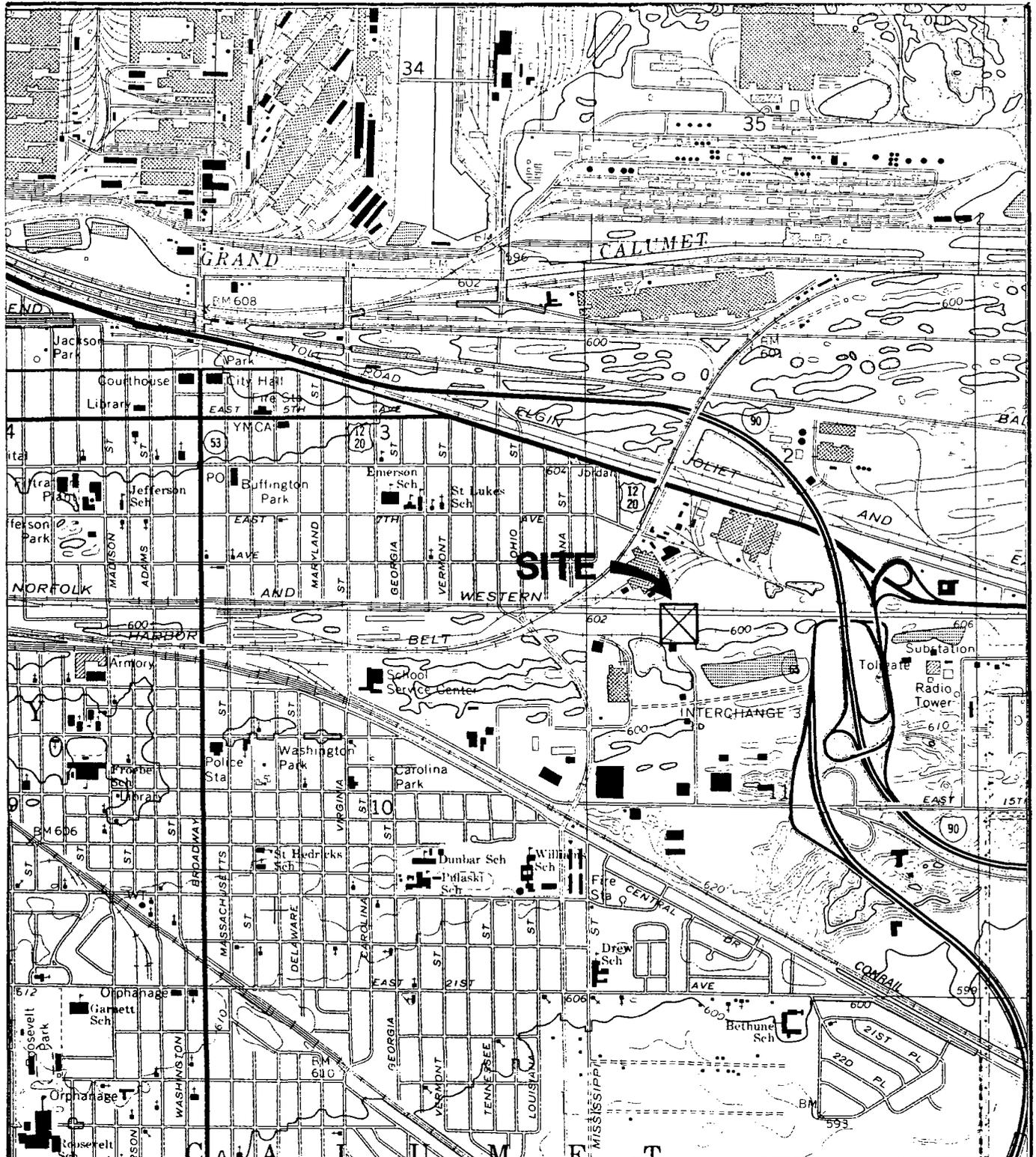


### Immediate Removal Action Check Sheet

	High	Moderate	Low
<b><u>Fire and Explosion Hazard</u></b>			
Flammable Materials <u>Phosphorus is ignitable</u>		X ←	X
Explosives <u>N/A</u>			
Incompatible Chemicals <u>Unknown</u>			X
<b><u>Direct Contact with Acutely Toxic Chemicals</u></b>			
Site Security <u>Fenced, gate locked at night</u>			X
Leaking Drums or Tanks <u>None</u>			X
Open Lagoons or pits <u>Open catch basin</u>		X	
Materials on Surface <u>contaminated soil/sed.</u>		X	
Proximity of Population <u>1850'</u>			X
Evidence of Casual Site Use <u>None</u>			X
<b><u>Contaminated Water Supply</u></b>			
Exceeds 10 Day Snarl <u>N/A</u>			
Gross Taste or Odors <u>N/A</u>			
Alternate Water Available <u>Yes</u>			X
Potential Contamination <u>No</u>			X
Is the site abandoned or <b>active?</b>			
<u>Leased by Gary Products Corp.</u>			

Comments

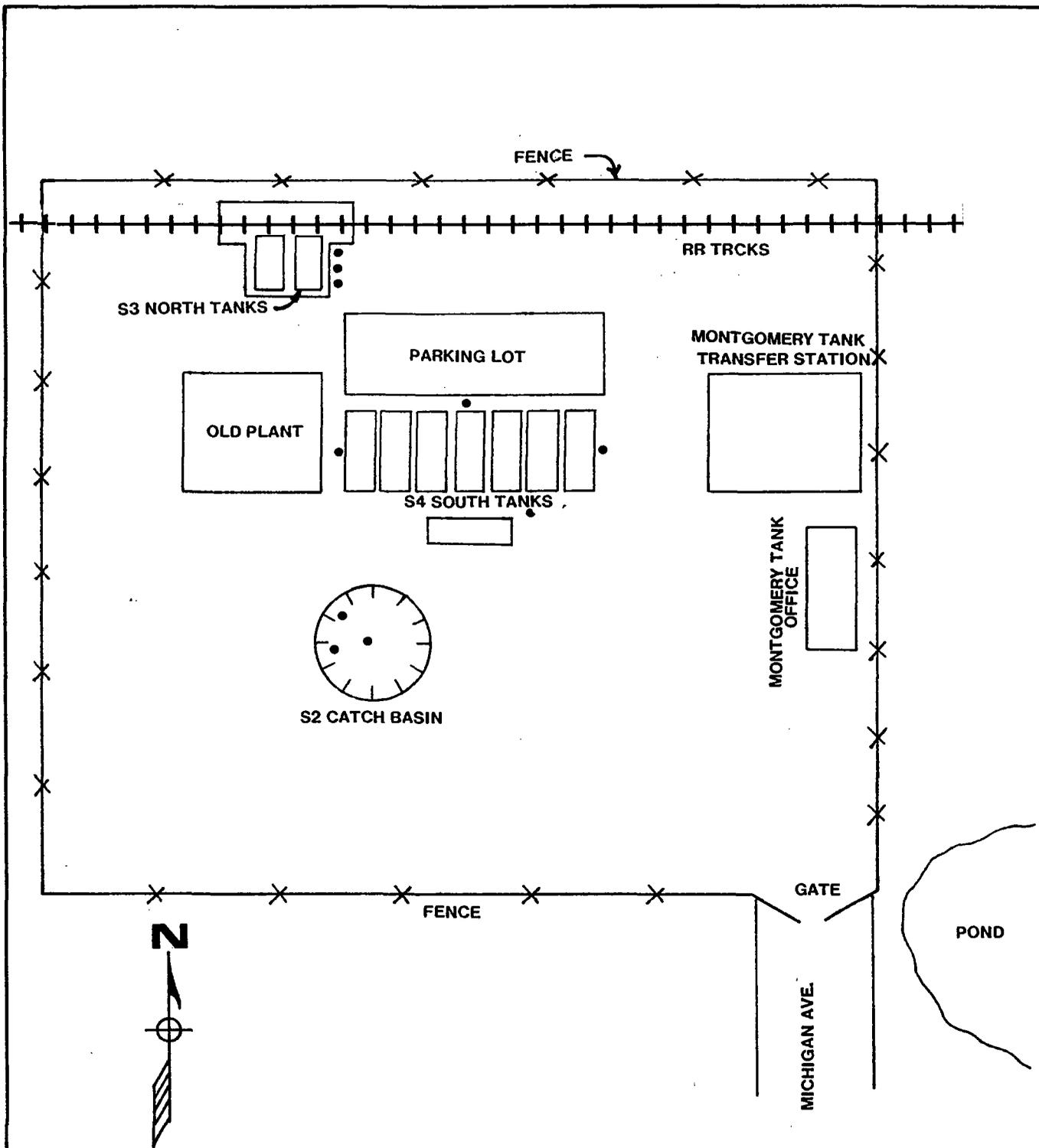




**ecology and environment, inc.**

111 WEST JACKSON BLVD., CHICAGO, ILLINOIS 60604, TEL. 312-663-9415

TITLE	<b>SITE LOCATION MAP</b>	FIGURE #	<b>1</b>
SITE	<b>MOBIL CHEMICAL PHOSPHORUS DIV.</b>	SCALE	<b>1:24,000</b>
CITY	<b>GARY, INDIANA</b>	TDD #	<b>F05-8702-175</b>
SOURCE	<b>USGS</b>	DATE	<b>1968</b>
		REVISED	<b>1980</b>



• COMPOSITE  
SAMPLE LOCATIONS

<b>ecology and environment, inc.</b> 111 WEST JACKSON BLVD., CHICAGO, ILLINOIS 60604, TEL. 312-683-9415			
<b>TITLE</b> SAMPLE LOCATION MAP		<b>FIGURE #</b> 2	
<b>SITE</b> MOBIL CHEMICAL PHOSPHORUS DIV		<b>SCALE</b> NONE	
<b>CITY</b> GARY, INDIANA		<b>IDO #</b> F05-8702-175	
<b>SOURCE</b> FIT SITE INSPECTION		<b>DATE</b> 2/18/87	
		<b>REVISED</b> N/A	

DATE 2/18/87

TIME 1:18 A.M. P.M.

DIRECTION: N (NNE) NE ENE  
E ESE SE SSE  
S SSW SW WSW  
W WNW NW NNW

WEATHER clear, 37°F

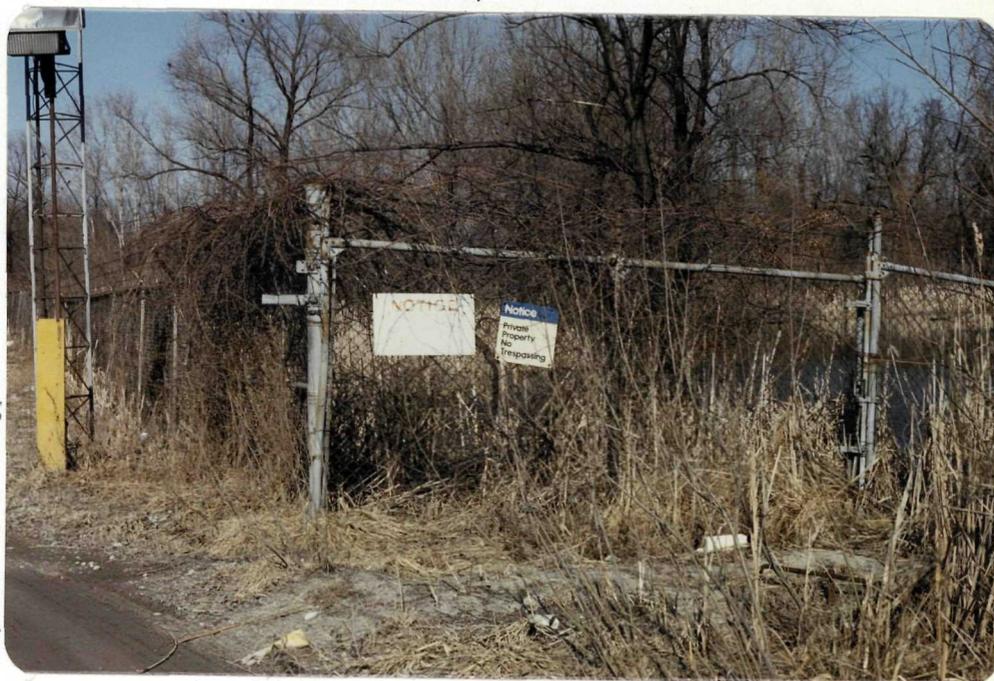
SITE Mobil Chemical-Phosphorus

TDD# F05-8702-172

PHOTOGRAPHED BY:

Julie Kaiser

SAMPLE ID# (if applicable)



DESCRIPTION:

DATE 2/18/87

TIME 1:18 A.M. P.M.

DIRECTION: (N) NNE NE ENE  
E ESE SE SSE  
S SSW SW WSW  
W WNW NW NNW

WEATHER clear, 37°F

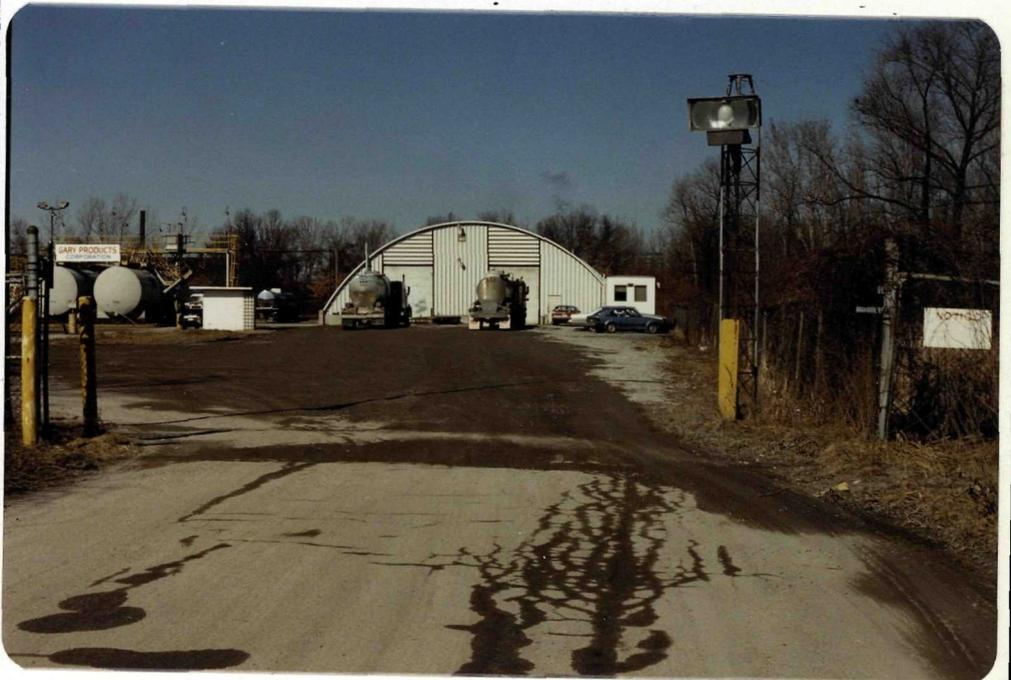
SITE Mobil Chemical-Phosphorus

TDD# F05-8702-172

PHOTOGRAPHED BY:

Julie Kaiser

SAMPLE ID# (if applicable)



DESCRIPTION:

DATE 2/18/87

TIME 12:15 A.M. (P.M.)

DIRECTION: N NNE NE ENE  
E ESE SE SSE  
S SSW SW WSW  
W WNW NW (ENE)

WEATHER clear, 37°F

SITE Mobil Chemical-Phosphorus

TDD# F05-8702-172

PHOTOGRAPHED BY:  
Julie Kaiser

SAMPLE ID# (if applicable)

DESCRIPTION:



DATE 2/18/87

TIME 12:25 A.M. (P.M.)

DIRECTION: N NNE NE (ENE)  
E ESE SE SSE  
S SSW SW WSW  
W WNW NW NNW

WEATHER clear, 37°F

SITE Mobil Chemical-Phosphorus

TDD# F05-8702-172

PHOTOGRAPHED BY:  
Julie Kaiser

SAMPLE ID# (if applicable)

S2 Lagoon

DESCRIPTION:



DATE 2/18/87.  
TIME 12:25 A.M. (P.M.)

DIRECTION: N NNE NE ENE  
E ESE SE SSE  
S SSW SW WSW  
W WNW NW NNW

WEATHER clear, 37°F

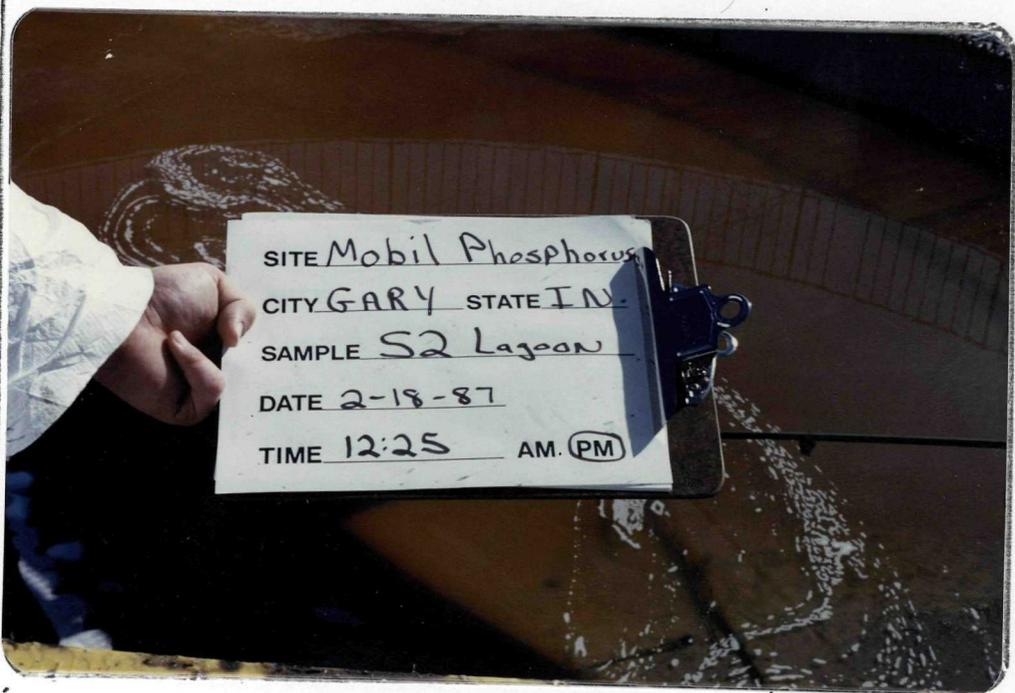
SITE Mobil Chemical-Phosphorus

ID# F05-8702-172

PHOTOGRAPHED BY:  
Julie Kaiser

SAMPLE ID# (if applicable)  
S2 Lagoon

DESCRIPTION:



DATE 2/18/87  
TIME 1:10 A.M. (P.M.)

DIRECTION: N NNE NE ENE  
E ESE SE SSE  
S SSW SW WSW  
W WNW NW NNW

WEATHER clear 37°F

SITE Mobil Chemical-Phosphorus

ID# F05-8702-172

PHOTOGRAPHED BY:  
Julie Kaiser

SAMPLE ID# (if applicable)

DESCRIPTION:



DATE 2/18/87

TIME 13:05 A.M. (P.M.)

DIRECTION: N NNE NE ENE  
E ESE SE SSE  
S SSW SW WSW  
W WNW NW NNW

WEATHER clear, 37°F

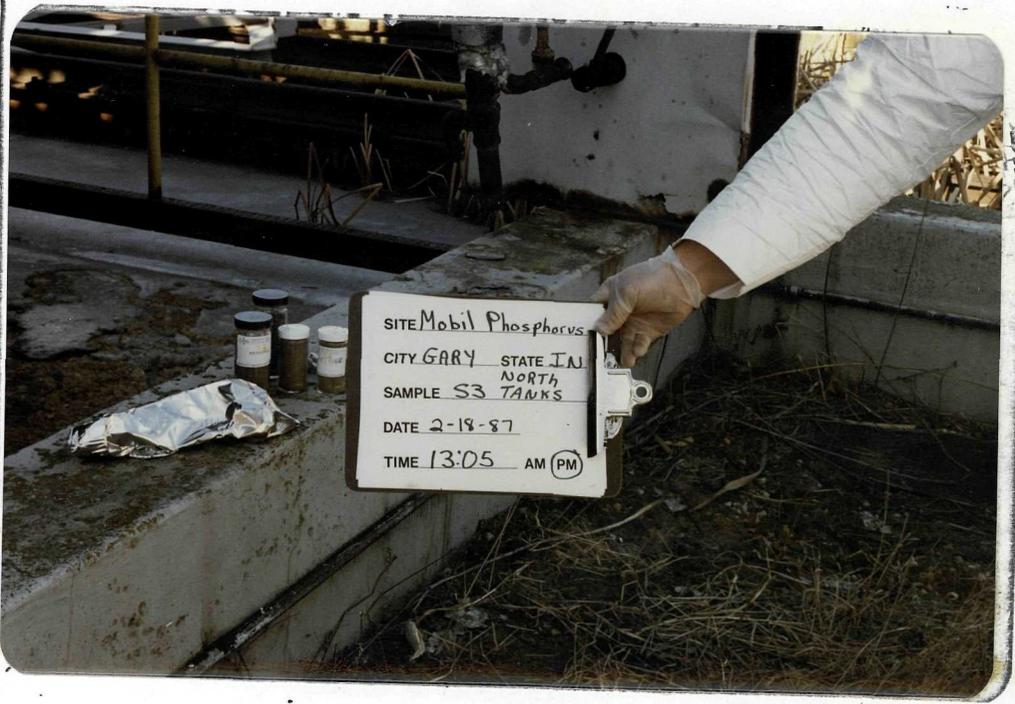
SITE Mobil Chemical-Phosphorus

ID# F05-8702-172

PHOTOGRAPHER BY:  
Julie Kaiser

SAMPLE ID# (if applicable)  
S3 North Tanks

DESCRIPTION:



DATE 2/18/87

TIME 13:06 A.M. (P.M.)

DIRECTION: N NNE NE ENE  
E ESE SE SSE  
S SSW SW WSW  
W WNW NW NNW

WEATHER clear, 37°F

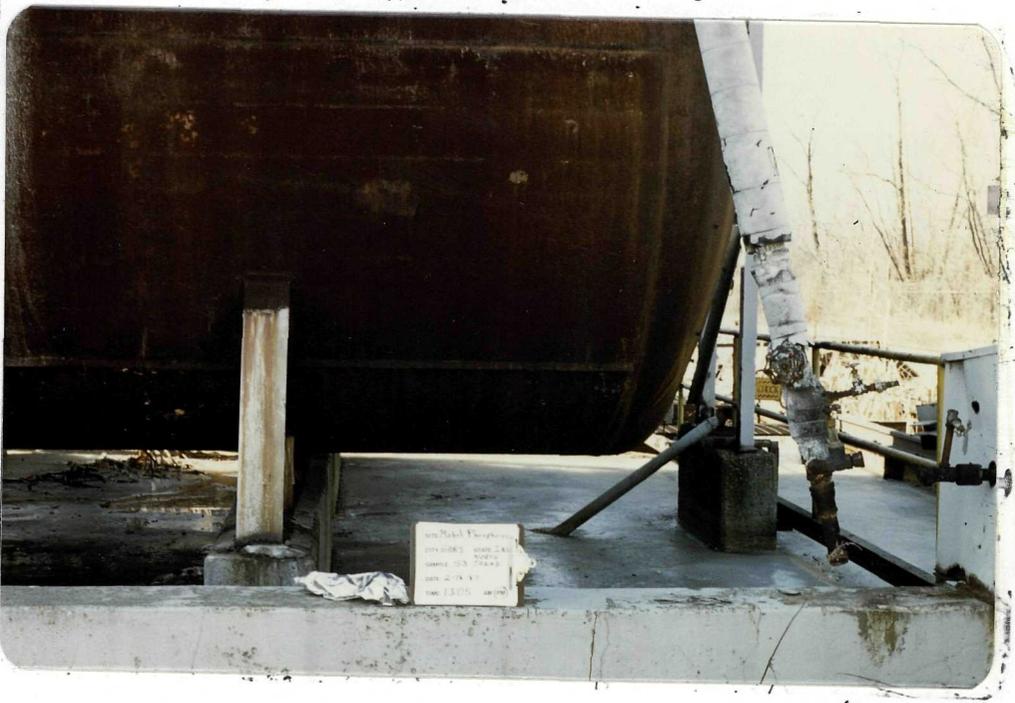
SITE mobil Chemical-Phosphorus

ID# F05-8702-172

PHOTOGRAPHER BY:  
Julie Kaiser

SAMPLE ID# (if applicable)  
S3 North Tanks

DESCRIPTION:



DATE 2/18/87

TIME 1:17 A.M. (P.M.)

DIRECTION: N NNE NE ENE  
E ESE SE SSE  
S SSW SW WSW  
W WNW NW (NNW)

WEATHER clear 37°F

SITE Mobil Chemical-Phosphorus

TDD# FOS-8702-172

PHOTOGRAPHED BY:  
Julie Kaiser

SAMPLE ID# (if applicable)



DESCRIPTION:

DATE 2/18/87

TIME 1:25 A.M. (P.M.)

DIRECTION: (N) NNE NE ENE  
E ESE SE SSE  
S SSW SW WSW  
W WNW NW NNW

WEATHER clear 37°F

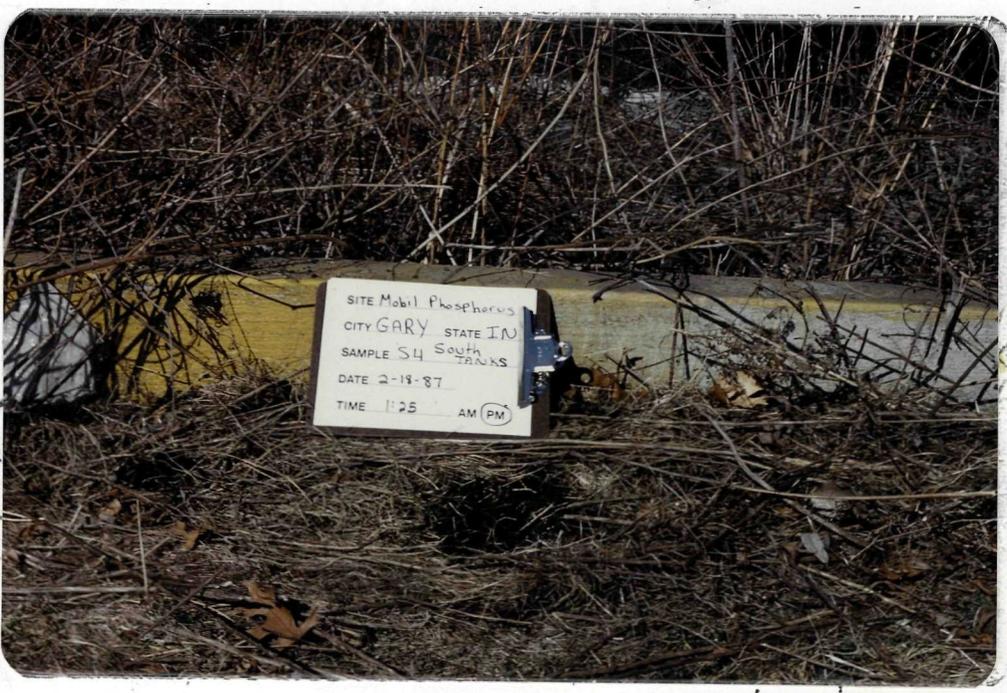
SITE Mobil Chemical-Phosphorus

TDD# FOS-8702-172

PHOTOGRAPHED BY:  
Julie Kaiser

SAMPLE ID# (if applicable)

SA South Tanks



DESCRIPTION:

DATE 2/18/87

TIME 1:25 A.M. (P.M.)

DIRECTION: (N) NNE NE ENE  
E ESE SE SSE  
S SSW SW WSW  
W WNW NW NNW

WEATHER clear 37°F

SITE Mobil Chemical - Prossphorus

ID# FOS-8702-172

PHOTOGRAPHED BY:  
Julie Kaiser

SAMPLE ID# (if applicable)

DESCRIPTION:



DATE 2/18/87

TIME 3:05 A.M. (P.M.)

DIRECTION: N NNE NE ENE  
E ESE SE SSE  
S SSW (SW) WSW  
W WNW NW NNW

WEATHER clear 37°F

SITE Mobil Chemical - Prossphorus

ID# FOS-8702-172

PHOTOGRAPHED BY:  
Julie Kaiser

SAMPLE ID# (if applicable)

DESCRIPTION:



DATE 2/18/87

TIME 3:00 A.M. (P.M.)

DIRECTION: N NNE NE ENE  
E ESE SE SSE  
S SSW (SW) WSW  
W WNW NW NNW

WEATHER clear 37° F

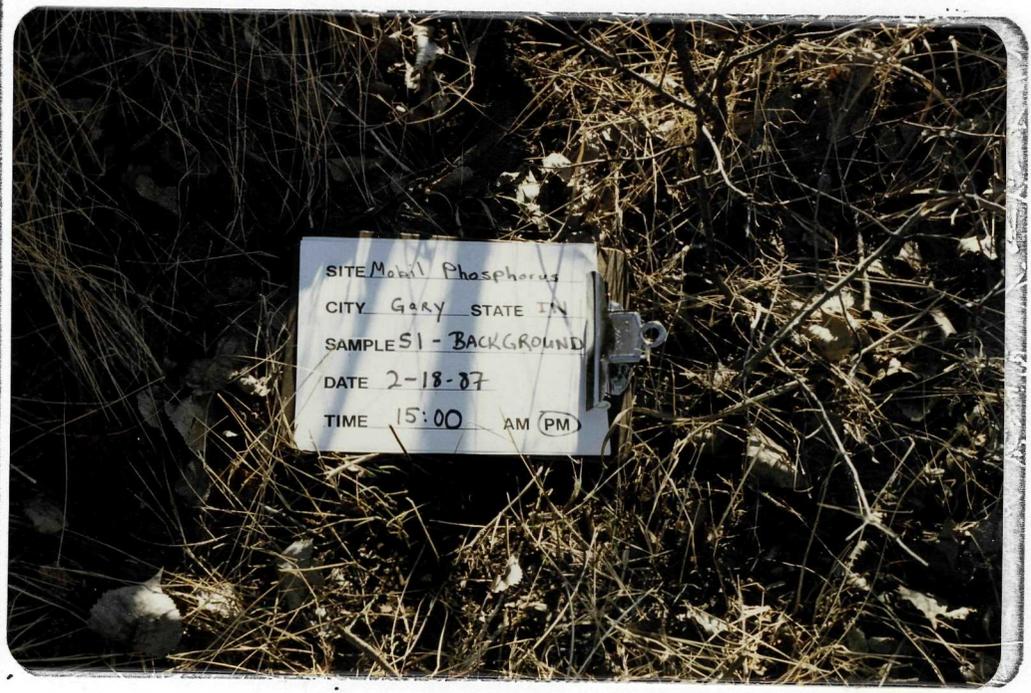
SITE Mobil Chemical-Phosphorus

DD# F05-8702-172

PHOTOGRAPHED BY:  
Julie Kaiser

SAMPLE ID# (if applicable)  
S1 Background

DESCRIPTION:



DATE

TIME A.M. P.M.

DIRECTION: N NNE NE ENE  
E ESE SE SSE  
S SSW SW WSW  
W WNW NW NNW

WEATHER

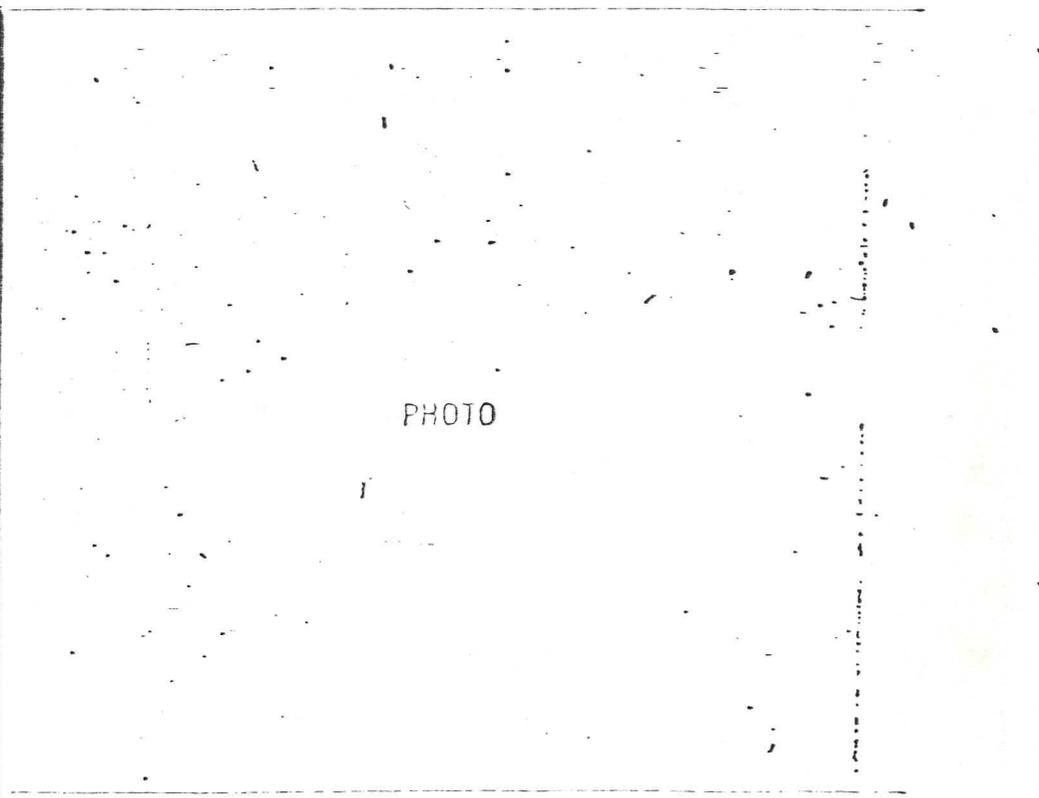
SITE

DD#

PHOTOGRAPHED BY:

SAMPLE ID# (if applicable)

DESCRIPTION:



PHOTO



COMPOUND	ITC			
	MEJ 375	MEJ 376	MEJ 377	MEI 948
	EJ 104	EJ 105	EJ 106	EJ 920
SAMPLE	S1	S2	S3	S4
	Background	catch basin	north tanks	south tanks
pentachlorophenol	130 J	55000		800 J
phenanthrene	180 J	5500 J		
anthracene				
di-n-butylphthalate	1300		1100 J	
fluoranthene	140 J		240 J	770 J
benzidine				
pyrene	120 J		250 J	540 J
butylbenzylphthalate				
3,3'-dichlorobenzidine				
benzo(a)anthracene	110 J			450 J
bis(2-ethylhexyl)phthalate		100 000	14000	
chrysene	120 J			540 J
di-n-octylphthalate				
benzo(b&k)fluoranthene	170 J			550 J
benzo(a)pyrene	100 J			560 J
indeno(1,2,3-cd)pyrene				
dibenzo(a,h)anthracene				
benzo(g,h,i)perylene				
alpha-BHC				
beta-BHC				
delta-BHC				
gamma-BHC(lindane)				
heptachlor				
aldrin				
heptachlor epoxide				
endosulfan I				
dieldrin				
4,4'-DDE				
endrin				
endosulfan II				
4,4'-DDD				
endrin aldehyde				
endosulfan sulfate				
4,4'-DDT				
methoxychlor				
endrin ketone				
chlordane				
toxaphene				
Aroclor-1016				
Aroclor-1221				
Aroclor-1232				
Aroclor-1242				
Aroclor-1248				
Aroclor-1254				
Aroclor-1260				
ELEMENT				
aluminum	1220	1020	1190	7840
antimony	[9.6] N			
arsenic	[2.7]			4.2
barium	[14]	[61]		100
beryllium				
cadmium		3.7		
calcium	2290	21,200	5810	60,000
chromium	7.0 *	81 *	6.7 *	19 *
cobalt			[4.3]	
copper	5.8	49	7.5	24
iron	5620	11400	6100	15400
lead	63 J	407 J	57 J	36 J
magnesium	[632]	4480	2820	26200
manganese	126	175	107	343
mercury		7.0	0.09	
nickel		8.1		
potassium	[328]	4070	[212]	1220
selenium				
silver				
sodium		[896]		[237]
thallium				
tin				
vanadium	[6.1]		[5.8]	12
zinc	98 *	407 *	49 *	96 *
cyanide CHECK IF ANALYZED (✓)				
TENTATIVELY IDENTIFIED ORGANICS				
2-Methyl propane		88 000 J		

COMPOUND	ITC	MEJ 375	MEJ 376	MEJ 377	MEJ 948
	OTC	EJ 104	EJ 105	EJ 106	EJ 920
	SAMPLE	S1 Background	S2 Catch basin	S3 north tanks	S4 South tanks
chloromethane					
bromomethane					
vinyl chloride					
chloroethane					
methylene chloride		9 B	13000 B	7 B	10 B
acetone		17 B	5600 JB	11 B	23 B
carbon disulfide					
1,1-dichloroethene			4100 J		
1,1-dichloroethane			9800 J		
trans-1,2,-dichloroethene			3000 J		
chloroform					
1,2-dichloroethane					
2-butanone			5100 JB		
1,1,1-trichloroethane			1 J		
carbon tetrachloride					
vinyl acetate					
bromodichloromethane					
1,1,2,2-tetrachloroethane					
1,2-dichloropropane					
trans-1,3-dichloropropene					
trichloroethene			5900 J		
dibromochloromethane					
1,1,2-trichloroethane					
benzene		41	7200 J	3 J	4 J
cis-1,3-dichloropropene					
2-chloroethylvinylether					
bromoform			3 JB		
2-hexanone					
4-methyl-2-pentanone					
tetrachloroethene			4600 J		
toluene		16 B	15000 B	4 JB	5 JB
chlorobenzene					
ethylbenzene		8 B	33000 B	3 JB	3 JB
styrene					
total xylenes		23 B	160000 B	12 B	11 B
N-nitrosodimethylamine					
phenol					
aniline					
bis(2-chloroethyl)ether					
2-chlorophenol					
1,3-dichlorobenzene					
1,4-dichlorobenzene					
benzyl alcohol					
1,2-dichlorobenzene			11000		
2-methylphenol					
bis(2-chloroisopropyl)ether					
4-methylphenol					
N-nitroso-di-n-propylamine					
hexachloroethane					
nitrobenzene					
isophrone				830 J	
2-nitrophenol					
2,4-dimethylphenol					
benzoic acid					
bis(2-chloroethoxy)methane					
2,4-dichlorophenol					
1,2,4-trichlorobenzene					
naphthalene			20000		
4-chloroaniline					
hexachlorobutadiene					
4-chloro-3-methylphenol					
2-methylnaphthalene		62 J	10000		
hexachlorocyclopentadiene					
2,4,6-trichlorophenol					
2,4,5-trichlorophenol					
2-chloronaphthalene					
2-nitroaniline					
dimethyl phthalate					
acenaphthylene					
3-nitroaniline			30000		
acenaphthene					
2,4-dinitrophenol					
4-nitrophenol					
dibenzofuran					
2,4-dinitrotoluene					
2,6-dinitrotoluene					
diethylphthalate					
4-chlorophenyl-phenylether					
fluorene					
4-nitroaniline					
4,6-dinitro-2-methylphenol					
N-nitrosodiphenylamine		140 J	20000		
4-bromophenyl-phenylether					
hexachlorobenzene					



# ecology and environment, inc.

111 WEST JACKSON BLVD., CHICAGO, ILLINOIS 60604, TEL. 312-663-9415

International Specialists in the Environment

Date Received for Review: 3/27/87 Date Review Completed: \_\_\_\_\_

TO: Julie Kaiser

FROM: Zena Gold-Kaufman ZK

SUBJECT: MOBIL CHEM  
F5-8012-012  
FINO11051

Sample Description: Case # 6855/2726E

Project Data Status: \*incomplete\* awaiting organics  
(4)

FIT Date Review Findings:

- All Se, Ag, and Cyanide data are estimated (J)
- All Pb data is estimated (J)
- All Sb is unusable.
- All Chromium data is estimated

Additional Comments:

\$ none

Book No. 6

Page No. 12



UNITED STATES ENVIRONMENTAL PROTECTION AGENCY  
REGION V

DATE: 3-24-87

RECEIVED MAR 26 1987

SUBJECT: Review of Region V CLP Data  
Received for Review on 3/4/87

FROM: Curtis Ross, Director  
Central Regional Laboratory *Jay Thecker*

TO: Data User: *Tit*

We have reviewed the data for the following case(s).

SITE NAME *Mobil Chem* SMO Case No. *68553A52726E*  
EPA Data Set No. *SF 3778* No. of Samples *4* D.U./Activity Numbers *Y9051C72100*

CRL No. *87FPO1559 - 87FPO1562*

SMO Traffic No. *MEJ375 - MEJ377 ; MEI948*

Contract Laboratory: *Century* Hours Required for Review: *12*

Following are our findings.

*This review covers four soil samples which were analyzed for metals and cyanide.*

*Spike recovery for Se (65%), Ag (44%) and CN (60%) are biased low and data should be construed as estimated (J).*

*Spike recovery for Pb is 22% and the sample results are quantitatively questionable - the results biased significantly low. The detection limit should be elevated.*

*Spike recovery for Sb is 11% and the possibility of false negatives exists. Data are unusable (R).*

*RPD for Cr is 37% and the results are estimated. Lab has done serial dilution on sample outside this core.*

- Data are acceptable for use. Sample MBI-780. *Ronald Piccini, 3/23/87*
- Data are acceptable for use with qualifications noted above.
- Data are preliminary - pending verification by Contractor Laboratory.
- Data are unacceptable.

cc: Dr. Alfred Haebeler/Joan Fisk/Cary Ward, EPA Support Services  
Ross K. Robeson, EMSL-Las Vegas  
Robert Pritchard, CLP/SMO

5F3778

REPORT NARRATIVE

RECEIVED MAR 26 1987

The following report contains the results for EPA case #6855. Total metals and cyanide analysis was performed on all four soil samples.

The analytical furnace spikes were performed by taking nine-tenths of the sample and adding one-tenth of a spiking solution. The calculation for analytical spike recovery is as follows:  $SSR-9/10 \text{ SR/SA} \times 100$ .

Please note that lead results were taken from the ICP analysis.

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U.S. EPA, CENTRAL REGIONAL LAB.  
536 S. CLARK STREET  
CHICAGO, ILLINOIS 60605

DATA REVIEW SHEET

WATER SAMPLE SPK. \_\_\_\_\_  
 WATER SAMPLE DUP. \_\_\_\_\_  
 SOIL SAMPLE SPK. MEJ-377  
 SOIL SAMPLE DUP. MEJ-377

MATRIX: soil  
 CONC.: low

SITE Metals Chem  
 LAB Century  
 REVIEWED BY P. Pichione

CASE # 6855  
 DATA SET # SF 3778  
 LAB Q.C.# 016  
 DATE: 3/23/87

Metals:	CALIBRATION		BLANKS		ICP INTER. SPIKE		DUPLICATE		LCS	linear range
	ICV	CCV	ICB	PREP	INT	FINAL	WATER	SOIL		
1. Aluminum	✓	✓	✓	✓	✓					✓
2. Antimony										
3. Arsenic										
4. Barium										
5. Beryllium										
6. Cadmium										
7. Calcium										37%
8. Chromium										
9. Cobalt										
10. Copper										
11. Iron										22%
12. Lead										
13. Magnesium										
14. Manganese										
15. Mercury										
16. Nickel										
17. Potassium										65%
18. Selenium										44%
19. Silver										
20. Sodium										
21. Thallium										
22. Vanadium										rpm 32%
23. Zinc										60%
24. Cyanide										
Other:										

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U.S. EPA Contract Laboratory Program  
Sample Management Office  
P.O. Box 818 - Alexandria, VA 22313  
703/557-2490 FTS: 8-557-2490

RECEIVED MAR 26 1007  
Date 3/3/07

COVER PAGE  
INORGANIC ANALYSES DATA PACKAGE

Lab Name CENTURY LABORATORIES

Case No. 6855

SOW No. 705

Q.C. Report No. 016

Lab Receipt Date 2/19/07

Sample Numbers

<u>EPA No.</u>	<u>Lab ID No.</u>	<u>EPA No.</u>	<u>Lab ID No.</u>
<u>MEJ-375</u>	<u>0016-01</u>		
<u>MEJ-376</u>	<u>0016-02</u>		
<u>MEJ-377</u>	<u>0016-03</u>		
<u>MEJ-944</u>	<u>0016-04</u>		

Comments:  
\_\_\_\_\_  
\_\_\_\_\_  
\_\_\_\_\_  
\_\_\_\_\_

ICP interelement and background corrections applied? Yes  No   
If yes, corrections applied before  or after  generation of raw data.

- Footnotes:
- NR - Not required by contract at this time
  - Form I:
  - Value - If the result is a value greater than or equal to the instrument detection limit but less than the contract-required detection limit, report the value in brackets (i.e., [10]). Indicate the analytical method used with P (for ICP), A (for Flame AA) or F (for Furnace AA).
  - U - Indicates element was analyzed for but not detected. Report with the instrument detection limit value (e.g., 10U).
  - E - Indicates a value estimated or not reported due to the presence of interference. Explanatory note included on cover page.
  - s - Indicates value determined by Method of Standard Addition.
  - N - Indicates spike sample recovery is not within control limits.
  - \* - Indicates duplicate analysis is not within control limits.
  - + - Indicates the correlation coefficient for method of standard addition is less than 0.995
  - M - Indicates duplicate injection results exceeded control limits.

Indicate method used: P for ICP; A for Flame AA and F for Furnace.

000003

U.S. EPA Contract Laboratory Program  
Sample Management Office  
P.O. Box 818 - Alexandria, VA 22313  
703/557-2490 FTS: 8-557-2490

EPA Sample No.  
MEJ-375

Date 3/3/87

INORGANIC ANALYSIS DATA SHEET

LAB NAME Century Laboratories, Inc. CASE NO. 6135  
SOW NO. 785  
LAB SAMPLE ID. NO. D016-01 QC REPORT NO. 016

Elements Identified and Measured

Concentration: Low  Medium   
Matrix: Water  Soil  Sludge  Other

ug/L or ng/kg dry weight (Circle One)

- |                     |                   |                                 |                  |
|---------------------|-------------------|---------------------------------|------------------|
| 1. <u>Aluminum</u>  | <u>1220 P</u>     | 13. <u>Magnesium</u>            | <u>1632 P</u>    |
| 2. <u>Antimony</u>  | <u>[9.6] PN</u>   | 14. <u>Manganese</u>            | <u>126 P</u>     |
| 3. <u>Arsenic</u>   | <u>[2.1] F</u>    | 15. <u>Mercury</u>              | <u>0.11 UCVA</u> |
| 4. <u>Barium</u>    | <u>[14] P</u>     | 16. <u>Nickel</u>               | <u>4.7 UP</u>    |
| 5. <u>Beryllium</u> | <u>1.2 UP</u>     | 17. <u>Potassium</u>            | <u>[324] P</u>   |
| 6. <u>Cadmium</u>   | <u>0.9 UP</u>     | 18. <u>Selenium</u>             | <u>0.46 UFN</u>  |
| 7. <u>Calcium</u>   | <u>2290 P</u>     | 19. <u>Silver</u>               | <u>2.34 PN</u>   |
| 8. <u>Chromium</u>  | <u>7.0 P*</u>     | 20. <u>Sodium</u>               | <u>117 UP</u>    |
| 9. <u>Cobalt</u>    | <u>2.3 UP</u>     | 21. <u>Thallium</u>             | <u>0.23 UF</u>   |
| 10. <u>Copper</u>   | <u>5.8 P</u>      | 22. <u>Vanadium</u>             | <u>[6.1] P</u>   |
| 11. <u>Iron</u>     | <u>5620 PA-16</u> | 23. <u>Zinc</u>                 | <u>98 P*</u>     |
| 12. <u>Lead</u>     | <u>63 PN</u>      | Percent Solids (%) <u>PL.32</u> |                  |
| <u>Cyanide</u>      | <u>0.73 UN</u>    |                                 |                  |

Footnotes: For reporting results to EPA, standard result qualifiers are used as defined on Cover Page. Additional flags or footnotes explaining results are encouraged. Definition of such flags must be explicit and contained on Cover Page, however.

Comments: MEDIUM BROWN SOIL

Lab Manager Kenneth Bond

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87 FPO/560

U.S. EPA Contract Laboratory Program  
Sample Management Office  
P.O. Box 818 - Alexandria, VA 22313  
703/557-2490 FTS: 8-557-2490

EPA Sample No.  
MEI-376

Date 3-3-87

INORGANIC ANALYSIS DATA SHEET

LAB NAME Century Laboratories, Inc.

CASE NO. 685

SOW NO. 785

LAB SAMPLE ID. NO. 0016-02

QC REPORT NO. 016

Elements Identified and Measured

Concentration: Low  Medium   
Matrix: Water  Soil  Sludge  Other

ug/L or mg/kg dry weight (Circle One)

1. <u>Aluminum</u> <u>1020P</u>	13. <u>Magnesium</u> <u>4480P</u>
2. <u>Antimony</u> <u>124PN</u>	14. <u>Manganese</u> <u>175P</u>
3. <u>Arsenic</u> <u>KS [LTI] LOW 2.2UF</u>	15. <u>Mercury</u> <u>7.0 CAAA</u>
4. <u>Barium</u> <u>[617P</u>	16. <u>Nickel</u> <u>8.1P</u>
5. <u>Beryllium</u> <u>2.0UP</u>	17. <u>Potassium</u> <u>4070P</u>
6. <u>Cadmium</u> <u>3.7P</u>	18. <u>Selenium</u> <u>0.88UPN</u>
7. <u>Calcium</u> <u>21,200P</u>	19. <u>Silver</u> <u>4.1UPN</u>
8. <u>Chromium</u> <u>81P*</u>	20. <u>Sodium</u> <u>[8967P</u>
9. <u>Cobalt</u> <u>4.1UP</u>	21. <u>Thallium</u> <u>0.44UF</u>
10. <u>Copper</u> <u>49P</u>	22. <u>Vanadium</u> <u>4.1UP</u>
11. <u>Iron</u> <u>11,400P</u>	23. <u>Zinc</u> <u>407P*</u>
12. <u>Lead</u> <u>407PN</u>	Percent Solids (%) <u>45.06</u>
<u>Cyanide</u> <u>0.44UN</u>	

Footnotes: For reporting results to EPA, standard result qualifiers are used as defined on Cover Page. Additional flags or footnotes explaining results are encouraged. Definition of such flags must be explicit and contained on Cover Page, however.

Comments: MEDIUM BROWN SOIL

Lab Manager Bennett Bond

00005

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U.S. EPA Contract Laboratory Program  
 Sample Management Office  
 P.O. Box 818 - Alexandria, VA 22313  
 703/557-2490 FTS: 8-557-2490

EPA Sample No.  
MEJ-377

Date 3/3/87

INORGANIC ANALYSIS DATA SHEET

LAB NAME Century Laboratories, Inc.

CASE NO. 6855

SOW NO. 785

LAB SAMPLE ID. NO. 0016-03

QC REPORT NO. 016

Elements Identified and Measured

Concentration: Low  Medium   
 Matrix: Water  Soil  Sludge  Other

ug/L or ug/kg dry weight (Circle One)

1. <u>Aluminum</u> <u>1190P</u>	13. <u>Magnesium</u> <u>2820P</u>
2. <u>Antimony</u> <u>8.74UPN</u>	14. <u>Manganese</u> <u>107P</u>
3. <u>Arsenic</u> <u>1.4UF</u>	15. <u>Mercury</u> <u>0.09 CVAA</u>
4. <u>Barium</u> <u>[14]P</u>	16. <u>Nickel</u> <u>5.8UP</u>
5. <u>Beryllium</u> <u>1.4UP</u>	17. <u>Potassium</u> <u>[212]P</u>
6. <u>Cadmium</u> <u>1.2UP</u>	18. <u>Selenium</u> <u>0.56UFN</u>
7. <u>Calcium</u> <u>5410P</u>	19. <u>Silver</u> <u>2.94UPN</u>
8. <u>Chromium</u> <u>6.7P*</u>	20. <u>Sodium</u> <u>145UP</u>
9. <u>Cobalt</u> <u>[4.3]P</u>	21. <u>Thallium</u> <u>0.21UF</u>
10. <u>Copper</u> <u>7.5P</u>	22. <u>Vanadium</u> <u>[5.4]P</u>
11. <u>Iron</u> <u>6100P</u>	23. <u>Zinc</u> <u>49P*</u>
12. <u>Lead</u> <u>57PN</u>	Percent Solids (%) <u>66.41</u>
<u>Cyanide</u> <u>0.29UN</u>	

Footnotes: For reporting results to EPA, standard result qualifiers are used as defined on Cover Page. Additional flags or footnotes explaining results are encouraged. Definition of such flags must be explicit and contained on Cover Page, however.

Comments: MEDIUM BROWN SOIL

Lab Manager Kenneth Bond

U.S. EPA Contract Laboratory Program  
Sample Management Office  
P.O. Box 818 - Alexandria, VA 22313  
703/557-2490 FTS: 6-557-2490

EPA Sample No.  
MEI-94X

Date 3/3/87

INORGANIC ANALYSIS DATA SHEET

LAB NAME Century Laboratories, Inc.

CASE NO. 6755

SOW NO. 785

LAB SAMPLE ID. NO. 0616-04

QC REPORT NO. 016

Elements Identified and Measured

Concentration: Low  Medium   
Matrix: Water  Soil  Sludge  Other

ug/L or ug/kg dry weight (Circle One)

1. <u>Aluminum</u> <u>7840P</u>	13. <u>Magnesium</u> <u>26,200P</u>
2. <u>Antimony</u> <u>7.34UPN</u>	14. <u>Manganese</u> <u>343P</u>
3. <u>Arsenic</u> <u>4.2F</u>	15. <u>Mercury</u> <u>0.09UCVAA</u>
4. <u>Barium</u> <u>100P</u>	16. <u>Nickel</u> <u>4.9UP</u>
5. <u>Beryllium</u> <u>1.2UP</u>	17. <u>Potassium</u> <u>1220P</u>
6. <u>Cadmium</u> <u>1.0UP</u>	18. <u>Selenium</u> <u>0.44UPN</u>
7. <u>Calcium</u> <u>60,000P</u>	19. <u>Silver</u> <u>2.4UPN</u>
8. <u>Chromium</u> <u>19P*</u>	20. <u>Sodium</u> <u>237P</u>
9. <u>Cobalt</u> <u>2.4UP</u>	21. <u>Tungsten</u> <u>0.22UF</u>
10. <u>Copper</u> <u>24P</u>	22. <u>Vanadium</u> <u>12P</u>
11. <u>Iron</u> <u>15,400P</u>	23. <u>Zinc</u> <u>96P*</u>
12. <u>Lead</u> <u>36PN</u>	Percent Solids (2) <u>85.00</u>
<u>Cyanide</u> <u>0.73UN</u>	

Footnotes: For reporting results to EPA, standard result qualifiers are used as defined on Cover Page. Additional flags or footnotes explaining results are encouraged. Definition of such flags must be explicit and contained on Cover Page, however.

Comments: Medium brown soil

Lab Manager Kenneth Bond

Q. C. Report No. 016

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BLANKS

LAB NAME CENTURY LABORAT

CASE NO. 6155

DATE 3-3-81

UNITS ug/l (SOIL P.B. ANALYSIS)

MATRIX SOIL

Compound	Initial Calibration Blank Value	Continuous Calibration				Preparation Blank	
		1	Blank Value		4	Matrix: WATER	Matrix: SOIL
			2	3		1	2
Metals:							
1. Aluminum	[307]	[35]	20u	20u	20u	[297]	[7.07]
2. Antimony	30u	[547]	30u	56	30u	54	<sup>12</sup> 6.0u
3. Arsenic	<del>50u</del> 3.0u	<del>5.0u</del> <sup>3.0u</sup> 3.0u	<del>5.0u</del> <sup>3.0u</sup> 3.0u	<del>5.0u</del> <sup>3.0u</sup> 3.0u	<del>5.0u</del> <sup>3.0u</sup> 3.0u	<del>5.0u</del> <sup>2.0u</sup> 2.0u	<del>1.0u</del> <sup>0.4u</sup> 0.4u
4. Barium	10u	10u	10u	10u	10u	10u	2.0u
5. Beryllium	5.0u	5.0u	5.0u	5.0u	5.0u	5.0u	1.0u
6. Cadmium	4.0u	4.0u	4.0u	4.0u	4.0u	4.0u	0.8u
7. Calcium	100u	100u	100u	100u	100u	[724]	[66]
8. Chromium	10u	10u	10u	10u	10u	10u	2.0u
9. Cobalt	[117]	[15]	10u	10u	10u	10u	[3.87]
10. Copper	6.0u	6.0u	6.0u	6.0u	6.0u	6.0u	1.2u
11. Iron	80u	80u	80u	80u	80u	80u	16u
12. Lead	35uP	35uP	35uP	35uP	35uP	35uP	7.0uP
13. Magnesium	100u	100u	100u	100u	100u	100u	20u
14. Manganese	5.0u	5.0u	5.0u	5.0u	5.0u	5.0u	1.0u
15. Mercury	<del>2.0u</del> 0.2u	<del>2.0u</del> <sup>0.2u</sup> 0.2u	<del>2.0u</del> <sup>0.2u</sup> 0.2u	<del>2.0u</del> <sup>0.2u</sup> 0.2u	<del>2.0u</del> <sup>0.2u</sup> 0.2u	<del>2.0u</del> <sup>0.2u</sup> 0.2u	<del>0.4u</del> <sup>0.1u</sup> 0.1u
16. Nickel	20u	20u	20u	20u	20u	20u	4.0u
17. Potassium	500u	500u	500u	500u	500u	500u	100u
18. Selenium	2.0u	2.0u	2.0u	2.0u		2.0u	0.4u
19. Silver	10u	10u	10u	10u	10u	10u	2.0u
20. Sodium	500u	500u	500u	500u	500u	500u	100u
21. Thallium	1.0u	1.0u	1.0u	1.0u	1.0u	1.0u	0.2u
22. Vanadium	10u	10u	10u	10u	10u	10u	2.0u
23. Zinc	6.0u	6.0u	6.0u	6.0u	6.0u	6.0u	[1.6]
Other:							
Cyanide	10u	10u	10u	10u	10u	10u	0.2u

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Form III-2

Q. C. Report No. 016

BLANKS

LAB NAME CENTURY LABORATORIES

CASE NO. 685

DATE 3-3-71

UNITS ug/l

Matrix SOIL

Compound	Initial Calibration Blank Value	Continuing Calibration Blank Value				Preparation Blank Matrix: Matrix:	
		1	2	3	4	1	2
Metals:							
1. Aluminum		20u					
2. Antimony		30u					
3. Arsenic		20u	3.0	3.0u			
4. Barium		10u					
5. Beryllium		5.0u					
6. Cadmium		4.0u					
7. Calcium		100u					
8. Chromium		10u					
9. Cobalt		10u					
10. Copper		6.0u					
11. Iron		50u					
12. Lead		35uP					
13. Magnesium		100u					
14. Manganese		5.0u					
15. Mercury							
16. Nickel		20u					
17. Potassium		500u					
18. Selenium							
19. Silver		10u					
20. Sodium		500u					
21. Thallium		1.0u	1.0u	1.0u			
22. Vanadium		10u					
23. Zinc		6.0u					
Other:							
Cyanide							

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FORM V

Q. C. Report No. 016  
SPIKE SAMPLE RECOVERY

LAB NAME CENTURY LABORATORIES  
DATE 3-3-87

CASE NO. 6855  
EPA Sample No. MES-377  
Lab Sample ID No. 0016-03  
Units ML/KL

Matrix SOIL

Compound	Control Limit ZR	Spiked Sample Result (SSR)	Sample Result (SR)	Spiked Added (SA)	ZR <sup>1</sup>
Metals:					
1. Aluminum	75-125	<del>820</del> NR	820	NR	*
2. Antimony	-	[117] <del>6.00</del>	6.00	100	11N
3. Arsenic	-	9.6	1.00	8.0	120
4. Barium	-	X 396	[9.8]	400	96
5. Beryllium	-	8.9	1.00	10	89
6. Cadmium	-	7.9	0.80	10	79
7. Calcium	-	NR	4000	NR	NR
8. Chromium	-	42	4.6	40	93
9. Cobalt	-	97	[3.0]	100	97
10. Copper	-	49	5.2	50	87
11. Iron	-	NR	4200	NR	NR
12. Lead	-	61	39	100	22N
13. Magnesium	-	NR	1940	NR	NR
14. Manganese	-	172	74	100	98
15. Mercury	-	0.55	0.09	0.50	92
16. Nickel	-	95	4.00	100	95
17. Potassium	-	NR	[146]	NR	NR
18. Selenium	-	1.3	0.390	2.0	65N
19. Silver	-	X 4.4	2.00	10	44N
20. Sodium	-	NR	100.400	NR	NR
21. Thallium	-	10	0.190	10	100
22. Vanadium	-	89	[4.0]	100	89
23. Zinc	-	125	34	100	91
Other:					
Cyanide	-	1.2	0.200	2.0	60N

<sup>1</sup> ZR = [(SSR - SR)/SA] x 100

\* - out of control

NR - Not required

Comments:

000015

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Form VI

Q. C. Report No. 016

DUPLICATES

LAB NAME CENTURY LABORATORIES

CASE NO. 685

DATE 3-3-87

EPA Sample No. ME1-377

Lab Sample ID No. 0016-03

Units MG/KG

MATRIX SOIL

Compound	Control Limit <sup>1</sup>	Sample(S)	Duplicate(D)	RPD <sup>2</sup>
<b>Metals:</b>				
1. Aluminum		820	871	6.0
2. Antimony		6.0u	5.9u	NC
3. Arsenic		1.0u	[1.3]	NC
4. Barium		[9.8]	[117]	NC
5. Beryllium		1.0u	1.0u	NC
6. Cadmium		0.8u	1.1	NC
7. Calcium		4000	4158	3.9
8. Chromium	±2.0	4.6	6.7	37*
9. Cobalt		[3.0]	[2.2]	NC
10. Copper		5.2	4.7	10
11. Iron		4200	5148	20
12. Lead		39	47	19
13. Magnesium		1940	1900	2.1
14. Manganese		74	81	9.0
15. Mercury		0.09	0.1	10
16. Nickel		4.0u	4.0u	NC
17. Potassium		[146]	[154]	NC
18. Selenium		0.3tu	0.3tu	NC
19. Silver		2.0u	2.0u	NC
20. Sodium		10040u	100u	NC
21. Thallium		0.19u	0.19u	NC
22. Vanadium		[4.07]	[4.8]	NC
23. Zinc		34	47	32*
<b>Other:</b>				
Cyanide		0.2u	0.2u	NC

\* Out of Control

<sup>1</sup> To be added at a later date.

<sup>2</sup> RPD =  $[|S - D| / ((S + D) / 2)] \times 100$

NC - Non calculable RPD due to value(s) less than CRDL.

INORGANICS  
LABORATORY RESPONSE TO RESULTS OF CCS

Criterion	Comments
B1	ENCLOSED PLEASE FIND Form I for ME5-375 Corrected for flog
B2	ME1-941 Ca + Mg values do exceed screen page. Must be diluted
C	Form II-3 corrected for Ba
D	An Ammonium blank was reported because this case was analyzed along with other water samples. Repeating forms were filled out for all three cases per simultaneously.
E1	Form IV, Pb mean and standard deviation taken from 5 just mean. Values are 4340, 4340, 4330, 4310, 4230
E2	ICS was diluted and screen for Mg. Look at spec that Al, Ca + Fe also appear on diluted analysis. Diluted value now appears on corrected Form IV
K1	ICP instrument does not print date + time for standardization. Please note it has been filled in. This has never been a problem before. Why now? Is this a new requirement?
K2	Initial run of deep blk showed Lizz Pb when screen again Pb was O.K. Only Pb value should have been crossed out. See corrected data page.



5F377.8

COVER SHEET

LABORATORY RESPONSE TO RESULTS OF  
CONTRACT COMPLIANCE SCREENING (CCS)

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MAR 10 1987

US EPA CENTRAL REGIONAL LAB.  
535 S. CLARK STREET  
CHICAGO, ILLINOIS 60605

Response To: (check one)

Organics CCS

Inorganics CCS

Response materials sent to Organics CCS should be sent to the attention of Doris Ling, SMO.

Response materials sent to Inorganics CCS should be sent to the attention of Sa'ad Masri, SMO.

Laboratory Name CENTURY LABORATORIES

Response Date 3-17-87

EPA Contract No. 731F

Date Screening Results Received at Laboratory 3-11-87

Case No. 6785

Sample Nos.\*  
\_\_\_\_\_  
\_\_\_\_\_  
\_\_\_\_\_  
\_\_\_\_\_

\*Only list sample numbers that require reconciliation.

This form is used to identify materials sent in response to results of Contract Compliance Screening (CCS). A separate form must accompany the response for each Case.

Please indicate (on the attached continuation form) which fractions and/or which criteria correspond with your resubmission. Response materials sent to CCS should also be copied to the Region and to EMSL/LV, each with this blue Cover Sheet.

SF 3778

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CONTRACT COMPLIANCE SCREENING SUMMARY FOR INORGANIC B PAGE 1 OF 3

CASE: 6855 CONTRACT: 68-01-731B SAMPLES: 4 GCNUM: 016 DATE: 3/10/87

LAB NAME: CENTRY MAR 13 1987 REGION: 5 SCREENED BY: SR

SAMPLE NO.	A COVER PAGE	B DATA SHEETS	C DATA CALIBRATION SHEETS	D CONT. BLANK	E ICS	F MATRIX SPIKE	G DUP.	H IDL	I LCG	J MSA	K HOLD. TIME	L RAW DATA	M TR IDL	N RANGE
MEJ375-11	IRS	RS	"	"	"	"	"	"	"	"	"	"	"	"
MEJ376-11	"	"	"	"	"	"	"	"	"	"	"	"	"	"
MEJ377-11	IRS	EX	IRS	EX	IRS	SU	"	"	"	"	"	"	"	"
MEJ378-11	"	"	"	"	"	"	"	"	"	"	"	"	"	"
MEJ379-11	"	"	"	"	"	"	"	"	"	"	"	"	"	"
MEJ380-11	"	"	"	"	"	"	"	"	"	"	"	"	"	"
MEJ381-11	"	"	"	"	"	"	"	"	"	"	"	"	"	"
MEJ382-11	"	"	"	"	"	"	"	"	"	"	"	"	"	"
MEJ383-11	"	"	"	"	"	"	"	"	"	"	"	"	"	"
MEJ384-11	"	"	"	"	"	"	"	"	"	"	"	"	"	"
MEJ385-11	"	"	"	"	"	"	"	"	"	"	"	"	"	"
MEJ386-11	"	"	"	"	"	"	"	"	"	"	"	"	"	"
MEJ387-11	"	"	"	"	"	"	"	"	"	"	"	"	"	"
MEJ388-11	"	"	"	"	"	"	"	"	"	"	"	"	"	"
MEJ389-11	"	"	"	"	"	"	"	"	"	"	"	"	"	"
MEJ390-11	"	"	"	"	"	"	"	"	"	"	"	"	"	"
MEJ391-11	"	"	"	"	"	"	"	"	"	"	"	"	"	"
MEJ392-11	"	"	"	"	"	"	"	"	"	"	"	"	"	"
MEJ393-11	"	"	"	"	"	"	"	"	"	"	"	"	"	"
MEJ394-11	"	"	"	"	"	"	"	"	"	"	"	"	"	"
MEJ395-11	"	"	"	"	"	"	"	"	"	"	"	"	"	"
MEJ396-11	"	"	"	"	"	"	"	"	"	"	"	"	"	"
MEJ397-11	"	"	"	"	"	"	"	"	"	"	"	"	"	"
MEJ398-11	"	"	"	"	"	"	"	"	"	"	"	"	"	"
MEJ399-11	"	"	"	"	"	"	"	"	"	"	"	"	"	"
MEJ400-11	"	"	"	"	"	"	"	"	"	"	"	"	"	"

PROBLEM CODE: EX = EXPLAIN, RS = RESUBMIT, SU = SUBMIT, NC = NONCOMPLIANT

SAMPLE NO.	IC/M	RECD.	DATE	ANAL.	INIT.	CALIBRATION	BLANKS	ICB	ICS	LINEAR RANGE	LCS	SERIAL DILUTION	PROBLEMS
ME1948-1	ILS	19FEB87	12:29	"	11:00	ICV	CCB	ICB	ICS	10:57 11:04 11:16	11:09 11:19 11:21		(R1) MET 375: "N" flag is required on Rb, not on Fe. (R5)
MEJ375-1	ILS	19FEB87	12:11	"	11:02	CCV	CCB			11:29 11:33	11:27		(R2) MEI 948: Values above linear range. Ca - 244,710 ppb (upper limit 109,000 ppb) Mg - 106,906 ppb (upper limit 50,000 ppb) Dilution is required. (S4)
MEJ376-1	ILS	19FEB87	"	"	"	CCV	CCB			12:04 12:07	12:02		(C) Form II-3. %R for Ba = 10.2% # 203% (R5)
MEJ377-1	ILS	19FEB87	12:17	"	"								(D) Form III: Why was an aqueous prep. Hk reported? (EX)
MEJ377-8	ILS	19FEB87	12:25	"	"								(E) Form IV: "P" move and "standard deviation" based on at least 5 measurements for Pb is required on Form IV. (R5)
MEJ377-9	ILS	19FEB87	12:20	"	"	CCV	CCB			12:42 12:44	12:38		(E) ICS: Values above the linear range of Rb, Rf, Ca, and Mg. Dilute and rerun. (S4)
						CCV	CCB			13:13 13:15	13:11		(K) ICP standardization: Neither the date nor the time is on the data. (R5)
						CCV	CCB			13:38 13:40	13:28		(K2) Why was ICP data cancelled and rerun? prep. Hk - 11:14 CCB - 12:00 / 12:35 CCV - 12:40, 13:30 (EX)

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CONTRACT COMPLIANCE SCREENING FOR INORGANICS WORKSHEET B - AA AND CN PAGE 3 OF 3  
 CASE: 6855 CONTRACT: 68-01-7318 SAMPLES: 4 GCNUM: 016 DATE: 3/10/87  
 LAB NAME: CENTRY REGION: 5 SCREENED BY: SR

SAMPLE IC/M	DATE	as		hg		se		tl		cn		PROBLEMS
		F	IF	CV	F	F	F	C				
MEJ948-1	ILSI 19FEB87			3/6 I C P		3/1 581	3/21	3/28	3/25			(1) According to raw data (and using a dilution factor of 4) the % difference for Mg = 5.2%. Form IX has a value of 38 % diff for Mg. Please correct this value and resubmit the form. (RS) MK3, B4, H, E
MEJ375-1	ILSI 19FEB87											(K3) Value c/d off & ICP raw data. c/b (11:27) - Mg. (RS) MK2, B3, H, E
MEJ376-1	ILSI 19FEB87											(B3) Form I: Method of analysis is missing. MGT 375 - As. (RS) MK2, B4, H, E
MEJ377-1	ILSI 19FEB87											(B3) MEI 448: Descriptions & calibration, clarity, feature, and standards are required in the comments section on Form I. (RS) MK3, H, E
MEJ377-8	ILSI 19FEB87											(H) Circle with for LCS (RS) Also Se: 2.0 is the IDL for furnace not ICP (RS) MK3, H, E
MEJ377-9	ILSI 19FEB87											(I) As: Why was undetected raw data reported as <0.003, if the IDL on Form III is 0.005 ppm. (EX) MK3, H, E



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TABLE OF ICV'S & LCS'S

	<u>ICV</u>	<u>LCS</u>
ICP Analysis	ICV #1	WP 475 C3 (All Elements) WP 581 C2 Na, K, Mg, Ba, Ca WP 186
Furnace Analysis	ICV #2 (As, Se) ICV #4 (Pb, Tl)	WP 284 C1 (Pb, As, Se) WP 980 C2 (Tl)(2:100)
Hg Analysis	WS 378 C2	
Cn Analysis	WP 179 C2	

000002

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Form II - 1

Q. C. Report No. 016

INITIAL AND CONTINUING CALIBRATION VERIFICATION<sup>3</sup>

LAB NAME CENTURY LABORATORIES

CASE NO. 685

SOW NO. 75

UNITS ug/l

DATE 3-3-77

Compound	Initial Calib. <sup>1</sup>			Continuing Calibration <sup>2</sup>					
	True Value	Found	ZR	True Value	Found	ZR	Found	ZR	Method <sup>4</sup>
1. Aluminum	2160	2010	93	2000	1950	99	2050	104	P
2. Antimony	807	801	99	1000	1050	105	1040	104	P
3. Arsenic	41	44	107	50	49	98	45	90	F
4. Barium	2120	1940	91	2000	1890	94	1940	97	P
5. Beryllium	527	476	90	1000	965	96	966	97	P
6. Cadmium	455	446	91	200	185	92	188	94	P
7. Calcium	5150	46500	90	10,000	9350	94	9740	97	P
8. Chromium	529	482	91	200	195	97	199	99	P
9. Cobalt	496	489	99	1000	1000	100	1040	104	P
10. Copper	524	505	96	1000	948	95	910	98	P
11. Iron	2060	1920	93	10,000	9830	98	10,050	100	P
12. Lead	5160	4640	90	1,000	984	90	959	96	P
13. Magnesium	25,500	<del>24,000</del> 24,500	94	10,000	9850	99	10,100	101	P
14. Manganese	520	490	94	1,000	973	97	994	99	P
15. Mercury	0.90	1.07	119	1.0	1.17	117	1.20	120	CVA
16. Nickel	495	466	94	1000	999	100	1030	103	P
17. Potassium	51800	49500	93	10,000	9660	97	10,600	106	P
18. Selenium	43	47	109	25	24	96	24	96	F
19. Silver	495	496	100	1,000	1030	103	1000	100	P
20. Sodium	50300	47,600	95	10,000	9,000	90	10,030	100	P
21. Thallium	47	<sup>50</sup> 49	<sup>106</sup> 104	50	53	106	52	104	F
22. Vanadium	520	464	90	1,000	948	95	961	96	P
23. Zinc	8070	2850	94	1000	970	97	1010	101	P
Other:									
Cyanide	188	188	100	200	216	108	207	103	

<sup>1</sup> Initial Calibration Source SEE NARRATIVE <sup>2</sup> Continuing Calibration Source MID-LANSE STD

<sup>3</sup> Control Limits: Mercury and Tin 80-120; Other Metals 90-110; Cyanide 85-115

<sup>4</sup> Indicate Analytical Method Used: P - ICP; A - Flame AA; F - Furnace AA

000008

Form II-2

Q. C. Report No. 016

INITIAL AND CONTINUING CALIBRATION VERIFICATION<sup>3</sup>

LAB NAME CENTURY LABORATORIES

CASE NO. 655

SOW NO. 75

DATE 3-3-87

UNITS ug/L

Compound	Initial Calib. <sup>1</sup>			Continuing Calibration <sup>2</sup>					
	True Value	Found	ZR	True Value	Found	ZR	Found	ZR	Method <sup>4</sup>
Metals:									
1. Aluminum				2000	2150	107	2110	105	P
2. Antimony				1000	1100	110	1030	105	P
3. Arsenic	41	43	105	50	52	104	53	106	F
4. Barium				2000	2030	101	1970	99	P
5. Beryllium				1000	1000	100	986	99	P
6. Cadmium				200	191	95	186	93	P
7. Calcium				10,000	979	98	9710	97	P
8. Chromium				200	198	99	202	101	P
9. Cobalt				1000	1050	105	1040	104	P
10. Copper				1000	1020	102	999	100	P
1. Iron				10,000	10,300	103	10,100	101	P
12. Lead				1000	922	92	993	99	P
13. Magnesium				10,000	10,300	103	10,300	103	P
14. Manganese				1000	1020	102	1000	101	P
15. Mercury				1.0	0.96	96			CVA
16. Nickel				1000	1040	104	104	104	P
17. Potassium				10,000	11,000	110	10,500	105	P
18. Selenium				25	24	96			F
19. Silver				1,000	1010	101	1000	100	P
20. Sodium				10,000	10,600	106	10,100	101	P
21. Thallium				50	52	104	54	108	F
22. Vanadium				1,000	985	98	973	97	P
23. Zinc				1000	1030	103	1020	102	P
Other:									
Cyanide									

<sup>1</sup> Initial Calibration Source SEE NARRATIVE    <sup>2</sup> Continuing Calibration Source MID-RANGE STD.

<sup>3</sup> Control Limits: Mercury and Tin 80-120;    Other Metals 90-110; Cyanide 85-115

<sup>4</sup> Indicate Analytical Method Used: (P) - ICP; (A) - Flame AA; (F) - Furnace AA

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Q. C. Report No. 016

INITIAL AND CONTINUING CALIBRATION VERIFICATION<sup>3</sup>

LAB NAME CENTURY LABORATORIES

CASE NO. UHS

SOW NO. 75

DATE 3-5-87

UNITS ug/l

Compound	Initial Calib. <sup>1</sup>			Continuing Calibration <sup>2</sup>					
	True Value	Found	ZR	True Value	Found	ZR	Found	ZR	Method <sup>4</sup>
1. Aluminum				2000	2150	107			P
2. Antimony				1000	1100	110			P
3. Arsenic				50	48	96	46	92	F
4. Barium				2000	2030	101.5	45		P
5. Beryllium				1000	1010	101			P
6. Cadmium				200	195	97			P
7. Calcium				10,000	9850	99			P
8. Chromium				200	202	101			P
9. Cobalt				1000	1050	105			P
0. Copper				1000	1020	102			P
1. Iron				10,000	10400	104			P
2. Lead				1,000	1000	100			P
3. Magnesium				10,000	10,400	104			P
4. Manganese				1000	1030	103			P
5. Mercury									
6. Nickel				1000	1060	106			P
7. Potassium				10,000	10,600	106			P
8. Selenium									
9. Silver				1000	1000	100			P
0. Sodium				10,000	10,200	102			P
1. Thallium				50	50	100	52	104	F
2. Vanadium				1,000	999	100			P
3. Zinc				1,000	1040	104			P
4. Cyanide									

Initial Calibration Source \_\_\_\_\_ <sup>2</sup> Continuing Calibration Source MID-RANGE STD.

Control Limits: Mercury and Tin 80-120; Other Metals 90-110; Cyanide 85-115

Indicate Analytical Method Used: P - ICP; A - Flame AA; F - Furnace AA

000010

Q. C. Report No. 016

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INITIAL AND CONTINUING CALIBRATION VERIFICATION<sup>3</sup>

LABORATORY NAME CENTURY LABORATORIES

CASE NO. 688

SOW NO. 78

DATE 3-3-67

UNITS ug/l

Compound	Initial Calib. <sup>1</sup>			Continuing Calibration <sup>2</sup>					
	True Value	Found	ZR	True Value	Found	ZR	Found	ZR	Method <sup>4</sup>
Metals:									
1. Aluminum									
2. Antimony									
3. Arsenic									
4. Barium									
5. Beryllium									
6. Cadmium									
7. Calcium									
8. Chromium									
9. Cobalt									
10. Copper									
11. Iron									
12. Lead									
13. Magnesium									
14. Manganese									
15. Mercury									
16. Nickel									
17. Potassium									
18. Selenium									
19. Silver									
20. Sodium									
21. Thallium				50	49	98			F
22. Vanadium									
23. Zinc									
Other:									
Cyanide									

<sup>1</sup> Initial Calibration Source \_\_\_\_\_ <sup>2</sup> Continuing Calibration Source \_\_\_\_\_

<sup>3</sup> Control Limits: Mercury and Tin 80-120; Other Metals 90-110; Cyanide 85-115

<sup>4</sup> Indicate Analytical Method Used: P - ICP; A - Flame AA; F - Furnace AA

Form IV

Q. C. Report No. 016

ICP INTERFERENCE CHECK SAMPLE

LAB NAME CENTURY LABORATORIES

CASE NO. 6755

DATE 3-2-87

Check Sample I.D. ICS-XNB

Check Sample Source EMSL/LV

Units ug/l

Element	Control Limits <sup>1</sup>		True <sup>2</sup>	Initial		Final	
	Mean	Std. Dev.		Observed	ZR	Observed	ZR
Aluminum			503.000	496,000	99	572,000	114
Antimony				459,000	97	509,000	101
Arsenic				300u		300u	
Barium			472	425	90	423	90
Beryllium			456	437	96	436	96
Cadmium			964	797	83	812	84
Calcium			472,000	466,474	95	452,000	90
Chromium			915	893	91	921	93
Cobalt			471	423	88	419	88
Copper			509	507	99	516	102
Iron			191,000	179,500	90	145,000	92
Lead	4350	0.11	4630	4170	90	4290	93
Magnesium			497,000	474,000	95	534,000	107
Manganese			522	506	97	514	98
Nickel			913	833	91	853	93
Potassium				500u		500u	
Selenium							
Silver			916	931	94	939	95
Sodium				500u		500u	
Zinc							
Vanadium			471	423	90	427	91
Barium			948	866	91	895	94

Mean value based on n = 5

1 = value of EPA ICP Interference Check Sample or contractor standard.

CASE: EB55  
SA52726F

	1	2	3	4	5	6	7	8	9	10
	EJ	EJ	EJ	EJ						
SAMPLE DTE IIC	104	105	106	920						
SAMPLE	S1 Blghrd	S2 catch	S3 N. FRANKS	S4 S. FRANKS						
COMPOUND										
chloromethane										
bromomethane										
vinyl chloride										
chloroethane	9 B	18,000 B	7 B	10 B						
methylene chloride	77 B	17,200 B	1 B	23 B						
acetone										
carbon disulfide										
1,1-dichloroethane		4100 T								
1,1-dichloroethane		6800 T								
trans-1,2-dichloroethane		3000 T								
chloroform										
1,2-dichloroethane										
2-butanone		5100 BT								
1,1,1-trichloroethane			1 J							
carbon tetrachloride										
vinyl acetate										
bromodichloromethane										
1,1,2,2-tetrachloroethane										
1,1-dichloropropene										
trans-1,3-dichloropropene										
trichloroethane		5900 T								
1,1,2-trichloroethane										
benzene	(H)	7200 J	3 J	4 T						
cis-1,3-dichloropropene										
2-chloroethylvinylether				3 J B						
bromoform										
2-hexanone										
4-methyl-2-pentanone										
tetrachloroethane		4100 T								
toluene	16 B	15,000 B	4 T B	5 T B						
chlorobenzene										
ethylbenzene	8 B	33,000 B	3 J B	3 J B						
styrene										
total xylenes	23 B	11,000 B	12 B	11 B						
N-nitrosodimethylamine										
phenol										
aniline										
bis(2-chloroethyl) ether										
2-chlorophenol										
1,3-dichlorobenzene										
1,4-dichlorobenzene										
benzyl alcohol										
1,2-dichlorobenzene		71000								
2-methylphenol										
bis(2-chloroisopropyl) ether										
4-methylphenol										
N-nitroso-di-n-propylamine										
hexachloroethane										
nitrobenzene										
isophrone					820 J					
2-nitrophenol										
2,4-dimethylphenol										
benzoic acid										
bis(2-chloroethoxy)methane										
2,4-dichlorophenol										
1,2,4-trichlorobenzene										
naphthalene			20000							
4-chloroaniline										
hexachlorobutadiene										
4-chloro-3-methylphenol										
2-methylnaphthalene			6.2 J	10000						
hexachlorocyclopentadiene										
2,4,6-trichlorophenol										
2,4,5-trichlorophenol										
2-chloronaphthalene										
2-nitroaniline										
dimethyl phthalate										
acenaphthylene										
3-nitroaniline										
acenaphthene										
2,4-dinitrophenol										
4-nitrophenol										
dibenzofuran										
2,4-dinitrotoluene										
2,6-dinitrotoluene										
diethylphthalate										
4-chlorophenyl-phenylether										
fluorene										
4-nitroaniline										
4,6-dinitro-2-methylphenol										
N-nitrosodiphenylamine										
4-bromophenyl-phenylether										
hexachlorobenzene										

U = Use for HRS  
 J = Semi quantitative  
 = unmarked name  
 an artifact  
 TC

B = found in  
 blank  
 J = estimate

R = do not use  
 calibration  
 problem

CASE: 6055  
SAS 2746E

	1	2	3	4	5	6	7	8	9	10
SAMPLE OIC TIC	EJ	EJ	EJ	EJ						
	107	105	106	920						
	Bkgd.	S2 catch basin	S3 N. tanks	S4 S. tanks						
COMPOUND										
pentachlorophenol	130J	(15000)								
phenanthrene	180J	1800J								
anthracene										
di-n-butylphthalate	1900		1900J							
fluoranthene	140J		240J	990J						
benzidine				540J						
pyrene	120J		250J							
butylbenzylphthalate										
3,3'-dichlorobenzidine										
benzo(a)anthracene	110J	(10000)	(14000)	450J						
bis(2-ethylhexyl)phthalate										
chrysene	120J			540J						
di-n-octyl phthalate										
benzo(b,k)fluoranthene	170J			550J						
benzo(a)pyrene	100J			560J						
indeno(1,2,3-cd)pyrene										
dibenzo(a,h)anthracene										
benzo(g,h,i)perylene										
alpha-BHC										
beta-BHC										
delta-BHC										
gamma-BHC(lindane)										
heptachlor				BDL						
aldrin										
heptachlor epoxide										
endosulfan I										
dieldrin										
4,4'-DDE										
endrin										
endosulfan II										
4,4'-DDD										
endrin aldehyde										
endosulfan sulfate										
4,4'-DDT										
methoxychlor										
endrin ketone										
chlorodane										
toxaphene										
Aroclor-1016										
Aroclor-1221										
Aroclor-1232										
Aroclor-1242										
Aroclor-1248										
Aroclor-1254										
Aroclor-1260										
ELEMENT										
aluminum										
antimony										
arsenic										
barium										
beryllium										
cadmium										
calcium										
chromium										
cobalt										
copper										
iron										
lead										
magnesium										
manganese										
mercury										
nickel										
potassium										
selenium										
silver										
sodium										
thallium										
tin										
vanadium										
zinc										
cyanide										
CHECK IF ANALYZED ( )										
TENTATIVELY IDENTIFIED ORGANICS										



BDL  
Below  
detection  
Limit

○ = see for HRS  
□ = semi quantitative  
= unmarked means an artifact  
TC

87FP01560X

Sample Number

EJ-105

## Organics Analysis Data Sheet (Page 1)

Laboratory Name: ATI  
 Lab Sample ID No: EJ-105  
 Sample Matrix: SOIL  
 Data Release Authorized By: C Sites

Case No: 6855/SAS 2726E  
 QC Report No: —  
 Contract No: 6801-7014  
 Date Sample Received: 2-19-87

### Volatile Compounds

Concentration: Low Medium (Circle One)

Date Extracted/Prepared: 2-25-87

Date Analyzed: 2-25-87

Conc/Dil Factor:  $\left(\frac{10\text{mL}}{7.0\text{g}}\right)\left(\frac{3\text{mL}}{.01\text{mL}}\right)\left(\frac{1}{.562}\right) \frac{2200}{\text{PH}} 8.82$

Percent Moisture: (Not Decanted) 43.8

CAS Number		ug/l or ug/Kg (Circle One)
74-87-3	Chloromethane	22,000U
74-83-9	Bromomethane	↓
75-01-4	Vinyl Chloride	↓
75-00-3	Chloroethane	↓
75-09-2	Methylene Chloride	13,000B
67-64-1	Acetone	5,600JB
75-15-0	Carbon Disulfide	11,000U
75-35-4	1, 1-Dichloroethene	4,100J
75-34-3	1, 1-Dichloroethane	8,800J
156-60-5	Trans-1, 2-Dichloroethene	3,000J
67-66-3	Chloroform	11,000U
107-06-2	1, 2-Dichloroethane	↓
78-93-3	2-Butanone	5,100JB
71-55-6	1, 1, 1-Trichloroethane	20,000
56-23-5	Carbon Tetrachloride	11,000U
108-05-4	Vinyl Acetate	22,000U
75-27-4	Bromodichloromethane	11,000U

CAS Number		ug/l or ug/Kg (Circle One)
78-87-5	1, 2-Dichloropropane	11,000U
10061-02-6	Trans-1, 3-Dichloropropene	↓
79-01-6	Trichloroethene	5,900J
124-48-1	Dibromochloromethane	11,000U
79-00-5	1, 1, 2-Trichloroethane	↓
71-43-2	Benzene	7,200J
10061-01-5	cis-1, 3-Dichloropropene	11,000U
110-75-8	2-Chloroethylvinylether	22,000U
75-25-2	Bromoform	11,000U
108-10-1	4-Methyl-2-Pentanone	22,000U
591-78-6	2-Hexanone	↓
127-18-4	Tetrachloroethene	4,600J
79-34-5	1, 1, 2, 2-Tetrachloroethane	71,000U
108-88-3	Toluene	15,000B
108-90-7	Chlorobenzene	11,000U
100-41-4	Ethylbenzene	33,000B
100-42-5	Styrene	11,000U
	Total Xylenes	160,000B

#### Data Reporting Qualifiers

For reporting results to EPA, the following results qualifiers are used. Additional flags or footnotes explaining results are encouraged. However, the definition of each flag must be explicit.

<p><b>Value</b> If the result is a value greater than or equal to the detection limit, report the value</p> <p><b>U</b> Indicates compound was analyzed for but not detected. Report the minimum detection limit for the sample with the U (e.g., 10U) based on necessary concentration/dilution action. (This is not necessarily the instrument detection limit.) The footnote should read: U-Compound was analyzed for but not detected. The number is the minimum attainable detection limit for the sample</p> <p><b>J</b> Indicates an estimated value. This flag is used either when estimating a concentration for tentatively identified compounds where a 1:1 response is assumed or when the mass spectral data indicated the presence of a compound that meets the identification criteria but the result is less than the specified detection limit but greater than zero (e.g., 10J) If limit of detection is 10 ug/l and a concentration of 3 ug/l is calculated, report as 3J</p>	<p><b>C</b> This flag applies to pesticide parameters where the identification has been confirmed by GC/MS. Single component pesticides <math>\geq 10</math> ng/l in the final extract should be confirmed by GC/MS</p> <p><b>B</b> This flag is used when the analyte is found in the blank as well as a sample. It indicates possible/probable blank contamination and warns the data user to take appropriate action</p> <p><b>Other</b> Other specific flags and footnotes may be required to properly define the results. If used, they must be fully described and such description attached to the data summary report</p>
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Laboratory Name AT7  
 Case No 6855/893 2726 E

Sample Number  
EJ105

Organics Analysis Data Sheet  
 (Page 2)

Semivolatile Compounds

Concentration:  Low Medium (Circle One)  
 Date Extracted/Prepared 2/19/87  
 Date Analyzed: 3/3/87  
 Conc./Dil Factor: 5.26 ml / (2.505 x 10<sup>3</sup>) x 610  
 Percent Moisture (Decanted) \_\_\_\_\_

GPC Cleanup  Yes  No  
 Separatory Funnel Extraction  Yes  
 Continuous Liquid-Liquid Extraction  Yes

CAS Number		ug/l or ug/Kg (Circle One)
108-95-2	Phenol	6100U
111-44-4	bis(2-Chloroethyl)Ether	
95-57-8	2-Chlorophenol	
541-73-1	1,3-Dichlorobenzene	
106-46-7	1,4-Dichlorobenzene	
100-51-6	Benzyl Alcohol	
95-50-1	1,2-Dichlorobenzene	11,000
95-48-7	2-Methylnaphthalene	6100U
39638-32-9	bis(2-chloroisopropyl)Ether	
106-44-5	4-Methylnaphthalene	
621-64-7	N-Nitroso-Di-n-Propylamine	
67-72-1	Hexachlorocyclopentadiene	
98-95-3	Nitrobenzene	
78-59-1	Isophorone	
88-75-5	2-Nitrophenol	
105-67-9	2,4-Dimethylnaphthalene	
65-85-0	Benzoic Acid	30,000U
111-91-1	bis(2-Chloroethoxy)Methane	6100U
120-83-2	2,4-Dichlorophenol	
120-82-1	1,2,4-Trichlorobenzene	
91-20-3	Naphthalene	20,000
106-47-8	4-Chloroaniline	6100U
87-68-3	Hexachlorobutadiene	
59-50-7	4-Chloro-3-Methylnaphthalene	
91-57-6	2-Methylnaphthalene	10,000
77-47-4	Hexachlorocycloheptadiene	6100U
88-06-2	2,4,6-Trichlorophenol	
95-95-4	2,4,5-Trichlorophenol	30,000U
91-58-7	2-Chloronaphthalene	6100U
88-74-4	2-Nitroaniline	30,000U
131-11-3	Dimethyl Phthalate	6100U
208-96-8	Acenaphthylene	
99-09-2	3-Nitroaniline	30,000

CAS Number		ug/l or ug/Kg (Circle One)
83-32-9	Acenaphthene	6100U
51-28-5	2,4-Dinitrophenol	30,000
100-02-7	4-Nitrophenol	
132-64-9	Dibenzofuran	6100U
121-14-2	2,4-Dinitrotoluene	
606-20-2	2,5-Dinitrotoluene	
84-66-2	Diethylphthalate	
7005-72-3	4-Chloroethoxyphenyl ether	
86-73-7	Fluorene	
100-01-6	4-Nitroaniline	30,000U
534-52-1	4,6-Dinitro-2-Methylnaphthalene	
86-30-6	N-Nitrosodiphenylamine (1)	20,000
101-55-3	4-Bromoethoxyphenyl ether	6100U
118-74-1	Hexachlorobenzene	
87-86-5	Pentachlorophenol	55,000
85-01-8	Phenanthrene	13,500
120-12-7	Anthracene	6100U
84-74-2	Di-n-Butylphthalate	
206-44-0	Fluoranthene	
129-00-0	Pyrene	
85-68-7	Butybenzylphthalate	
91-94-1	3,3-Dichlorobenzidine	12,000U
56-55-3	Benzofluoranthene	6100U
117-81-7	bis(2-Ethylhexyl)Phthalate	100,000
218-01-9	Chrysene	6100U
117-84-0	Di-n-Octyl Phthalate	
205-99-2	Benzobifluoranthene	
207-08-9	Benzokifluoranthene	
50-32-8	Benzofluorene	
193-39-5	Indeno[1,2,3-cd]pyrene	
53-70-3	Dibenzofluoranthene	
191-24-2	Benzofluorene	

(1)-Cannot be separated from diphenylamine

Laboratory Name ATI  
 Case No 6855 / CAS 2726E

Sample Number  
1:10  
EJ105

Organics Analysis Data Sheet  
 (Page 3)

Pesticide/PCBs

Concentration: Low Medium (Circle One)  
 Date Extracted/Prepared: 19 FEB 87  
 Date Analyzed: 5 MARCH 87  
 Conc/Dil Factor: 0.1  
 Percent Moisture (decanted) 43.8

GPC Cleanup  Yes  No  
 Separatory Funnel Extraction  Yes  
 Continuous Liquid - Liquid Extraction  Yes

CAS Number		ug/l or ug/Kg (Circle One)
319-84-6	Alpha-BHC	140m
319-85-7	Beta-BHC	
319-86-8	Delta-BHC	
58-89-9	Gamma-BHC (Lindane)	
76-44-8	Heptachlor	
309-00-2	Aldrin	
1024-57-3	Heptachlor Epoxide	
959-98-8	Endosulfan I	↓
60-57-1	Dieldrin	280m
72-55-9	4, 4'-DDE	
72-20-8	Endrin	
33213-65-9	Endosulfan II	
72-54-8	4, 4'-DDD	
1031-07-8	Endosulfan Sulfate	
50-29-3	4, 4'-DDT	↓
72-43-5	Methoxychlor	1400m
53494-70-5	Endrin Ketone	280m
57-74-9	Chlordane	1400m
8001-35-2	Toxaphene	2800m
12674-11-2	Aroclor-1016	1400m
11104-28-2	Aroclor-1221	
11141-16-5	Aroclor-1232	
53469-21-9	Aroclor-1242	
12672-29-6	Aroclor-1248	↓
11097-69-1	Aroclor-1254	2800m
11096-82-5	Aroclor-1260	↓

$V_i$  = Volume of extract injected (ul)

$V_s$  = Volume of water extracted (ml)

$W_s$  = Weight of sample extracted (g)

$V_t$  = Volume of total extract (ul)

$V_s$  \_\_\_\_\_ or  $W_s$  30.5  $V_t$  20000  $V_i$  2

x

Laboratory Name: ATI  
Case No 6855/SAS 2726E

Sample Number  
**EJ-105**

Organics Analysis Data Sheet  
(Page 4)

Tentatively Identified Compounds

CAS Number	Compound Name	Fraction	RT or Scan Number	Estimated Concentration (ug/l or ug/kg)
1. <u>75-28-5</u>	<u>2-Methyl propanoic acid</u>	<u>VOA</u>	<u>216</u>	<u>88,000 J</u>
2. <u>-</u>	<u>an Acid</u>	<u>↓</u>	<u>613</u>	<u>4,400 J</u>
3.				
4.				
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30.				

Laboratory Name: ATI  
 Case No 6855/SAS 2726 E

Sample Number  
EJ105

Organics Analysis Data Sheet  
 (Page 4)

Tentatively Identified Compounds

CAS Number	Compound Name	Fraction	RT or Scan Number	Estimated Concentration (ug/l or ug/kg)
1. —	unknown $\approx$ C <sub>12</sub> hydrocarbon	PNA	1108	60,000 J
2. —	unk. branched hyd. $\approx$ C <sub>10</sub>		1207	78,000 J
3. —	unk. branched phenol		1313	340,000 J
4. —	unk. branched hyd. $\approx$ C <sub>13</sub>		1345	92,000 J
5. —	unk. branched hyd $\approx$ C <sub>17</sub>		1392	320,000 J
6. —	unk. branched phenol		1403	180,000 J
7. —	unk. branched phenol		1411	240,000 J
8. —	unknown		1416	200,000 J
9. —	unk. branched phenol		1442	130,000 J
10. —	unk. phthalate		1449	100,000 J
11. —	unknown		1532	200,000 J
12. —	unknown		1604	280,000 J
13. —	unknown		1614	360,000 J
14. —	unknown		1621	540,000 J
15. —	unk. branched alcohol		1629	220,000 J
16. —	unk. organic acid		1639	440,000 J
17. —	unknown		1653	480,000 J
18. —	unknown	N	1660	140,000 J
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87 FPD 15618

Sample Number  
EJ-106

Organics Analysis Data Sheet  
(Page 1)

Laboratory Name: ATI  
 Lab Sample ID No: EJ-106  
 Sample Matrix: SOIL  
 Data Release Authorized By: C. Sites

Case No: 6855/SAS 2726E  
 QC Report No: -  
 Contract No: 6801-7014  
 Date Sample Received: 2/19/87

Volatile Compounds

Concentration: Low Medium (Circle One)  
 Date Extracted/Prepared: 2-24-87  
 Date Analyzed: 2-24-87  
 Conc/Dil Factor: 5g/5g = 1.0 pH 7.71  
 Percent Moisture: (Not Decanted) 4.4

CAS Number		ug/l or ug/Kg (Circle One)
74-87-3	Chloromethane	10U
74-83-9	Bromomethane	
75-01-4	Vinyl Chloride	
75-00-3	Chloroethane	
75-09-2	Methylene Chloride	7B
67-64-1	Acetone	11B
75-15-0	Carbon Disulfide	5U
75-35-4	1, 1-Dichloroethane	
75-34-3	1, 1-Dichloroethane	
156-60-5	Trans-1, 2-Dichloroethene	
67-66-3	Chloroform	
107-06-2	1, 2-Dichloroethane	
78-93-3	2-Butanone	10U
71-55-6	1, 1, 1-Trichloroethane	1J
56-23-5	Carbon Tetrachloride	5U
108-05-4	Vinyl Acetate	10U
75-27-4	Bromodichloromethane	5U

CAS Number		ug/l or ug/Kg (Circle One)
78-87-5	1, 2-Dichloropropane	5U
10061-02-6	Trans-1, 3-Dichloropropene	
79-01-6	Trichloroethene	
124-48-1	Dibromochloromethane	
79-00-5	1, 1, 2-Trichloroethane	
71-43-2	Benzene	3J
10061-01-5	cis-1, 3-Dichloropropene	5U
110-75-8	2-Chloroethylvinylether	10U
75-25-2	Bromoform	3JB
108-10-1	4-Methyl-2-Pentanone	10U
591-78-6	2-Hexanone	
127-18-4	Tetrachloroethene	5U
79-34-5	1, 1, 2, 2-Tetrachloroethane	
108-88-3	Toluene	4JB
108-90-7	Chlorobenzene	5U
100-41-4	Ethylbenzene	3JB
100-42-5	Styrene	5U
	Total Xylenes	12B

Data Reporting Qualifiers

For reporting results to EPA, the following results qualifiers are used. Additional flags or footnotes explaining results are encouraged. However, the definition of each flag must be explicit.

- Value** If the result is a value greater than or equal to the detection limit, report the value
- U** Indicates compound was analyzed for but not detected. Report the minimum detection limit for the sample with the U (e.g., 10U) based on necessary concentration/dilution action. (This is not necessarily the instrument detection limit.) The footnote should read U-Compound was analyzed for but not detected. The number is the minimum attainable detection limit for the sample
- J** Indicates an estimated value. This flag is used either when estimating a concentration for tentatively identified compounds where a 1:1 response is assumed or when the mass spectral data indicated the presence of a compound that meets the identification criteria but the result is less than the specified detection limit but greater than zero (e.g., 10J). If limit of detection is 10 ug/l and a concentration of 3 ug/l is calculated, report as 3J
- C** This flag applies to pesticide parameters where the identification has been confirmed by GC/MS. Single component pesticides ≥ 10 ng / ul in the final extract should be confirmed by GC/MS
- B** This flag is used when the analyte is found in the blank as well as a sample. It indicates possible/probable blank contamination and warns the data user to take appropriate action
- Other** Other specific flags and footnotes may be required to properly define the results. If used, they must be fully described and such description attached to the data summary report

Laboratory Name ATI  
 Case No: 6855/SAS 2726 E

Sample Number  
EJ 106

**Organics Analysis Data Sheet**  
(Page 2)

**Semivolatile Compounds**

Concentration: (Low) Medium (Circle One)  
 Date Extracted / Prepared: 2/19/87  
 Date Analyzed: 2/25/87  
 Conc./Dil Factor:  $\frac{1 \text{ mL}}{2.19 \text{ (DSC)}} \times 10^3 \times 4 = \times 140$   
 Percent Moisture (Decanted): —

GPC Cleanup  Yes  No  
 Separatory Funnel Extraction  Yes  
 Continuous Liquid - Liquid Extraction  Yes

CAS Number		ug/l or ug/Kg (Circle One)
108-95-2	Phenol	1400 U
111-44-4	bis:2-Chloroethyl Ether	
95-57-8	2-Chlorophenol	
541-73-1	1,3-Dichlorobenzene	
106-46-7	1,4-Dichlorobenzene	
100-51-6	Benzyl Alcohol	
95-50-1	1,2-Dichlorobenzene	
95-48-7	2-Methoxyphenol	
39638-32-9	bis:2-chloroisopropyl Ether	
106-44-5	4-Methoxyphenol	
621-64-7	N-Nitroso-Di-n-Propylamine	
67-72-1	Hexachlorobenzene	
98-95-3	Nitrobenzene	
78-59-1	Isophorone	
88-75-5	2-Nitrophenol	
105-67-9	2,4-Dimethoxyphenol	71400 U
65-85-0	Benzoic Acid	7000 U
111-91-1	bis:2-Chloroethoxyethane	1400 U
120-83-2	2,4-Dichlorophenol	
120-82-1	1,2,4-Trichlorobenzene	
91-20-3	Naphthalene	
106-47-8	4-Chloroaniline	
87-68-3	Hexachlorobutadiene	
59-50-7	4-Chloro-3-Methoxyphenol	
91-57-6	2-Methylnaphthalene	
77-47-4	Hexachlorocyclopentadiene	
88-06-2	2,4,6-Trichlorophenol	1400 U
95-95-4	2,4,5-Trichlorophenol	7000 U
91-58-7	2-Chloronaphthalene	1400 U
88-74-4	2-Nitroaniline	7000 U
131-11-3	Dimethyl Phthalate	1400 U
208-96-8	Acenaphthylene	
99-09-2	3-Nitroaniline	7000 U

CAS Number		ug/l or ug/Kg (Circle One)
83-32-9	Acenaphthene	1400 U
51-28-5	2,4-Dinitrophenol	7000 U
100-02-7	4-Nitrophenol	
132-64-9	Dibenzofuran	1400 U
121-14-2	2,4-Dinitrotoluene	
606-20-2	2,6-Dinitrotoluene	
84-66-2	Diethylphthalate	
7005-72-3	4-Chlorophenylphenyl ether	
86-73-7	Fluorene	1400 U
100-01-6	4-Nitroaniline	7000 U
534-52-1	4,6-Dinitro-2-Methoxyphenol	
86-30-6	N-Nitrosodiphenylamine (1)	1400 U
101-55-3	4-Bromophenylphenyl ether	
118-74-1	Hexachlorobenzene	
87-86-5	Pentachlorophenol	7000 U
85-01-8	Phenanthrene	1400 U
120-12-7	Anthracene	
84-74-2	Di-n-Butylphthalate	1100 J
206-44-0	Fluoranthene	240 J
129-00-0	Pyrene	250 J
85-68-7	Butylbenzylphthalate	1400 U
91-94-1	3,3'-Dichlorobenzidine	2800 U
56-55-3	Benzofluoranthene	1400 U
117-81-7	bis:2-Ethylhexylphthalate	14000 U
218-01-9	Chrysene	1400 U
117-84-0	Di-n-Octyl Phthalate	
205-99-2	Benzodifluoranthene	
207-08-9	Benzokifluoranthene	
50-32-8	Benzofluoranthene	
193-39-5	Indeno[1,2,3-cd]pyrene	
53-70-3	Dibenzofluoranthene	
191-24-2	Benzofluoranthene	

(1)-Cannot be separated from diphenylamine

Laboratory Name ATI  
 Case No 6855/SAS 2726E

Sample Number  
7:5  
EJ 106

Organics Analysis Data Sheet  
 (Page 3)

Pesticide/PCBs

Concentration: Low Medium (Circle One)  
 Date Extracted/Prepared: 19 FEB 87  
 Date Analyzed: 4 MARCH 87  
 Conc/Dil Factor: 0.2  
 Percent Moisture (decanted) 4.44

GPC Cleanup  Yes  No  
 Separatory Funnel Extraction  Yes  
 Continuous Liquid - Liquid Extraction  Yes

CAS Number		ug/l or ug/Kg (Circle One)
319-84-6	Alpha-BHC	42m
319-85-7	Beta-BHC	
319-86-8	Delta-BHC	
58-89-9	Gamma-BHC (Lindane)	
76-44-8	Heptachlor	
309-00-2	Aldrin	
1024-57-3	Heptachlor Epoxide	
959-98-8	Endosulfan I	↓
60-57-1	Dieldrin	84m
72-55-9	4,4'-DDE	
72-20-8	Endrin	
33213-65-9	Endosulfan II	
72-54-8	4,4'-DDD	
1031-07-8	Endosulfan Sulfate	
50-29-3	4,4'-DDT	↓
72-43-5	Methoxychlor	420m
53494-70-5	Endrin Ketone	84m
57-74-9	Chlordane	420m
8001-35-2	Toxaphene	840m
12674-11-2	Aroclor-1016	420m
11104-28-2	Aroclor-1221	
11141-16-5	Aroclor-1232	
53469-21-9	Aroclor-1242	
12672-29-6	Aroclor-1248	↓
11097-69-1	Aroclor-1254	840m
11096-82-5	Aroclor-1260	↓

$V_i$  = Volume of extract injected (ul)  
 $V_s$  = Volume of water extracted (ml)  
 $W_s$  = Weight of sample extracted (g)  
 $V_t$  = Volume of total extract (ul)

$V_s$  \_\_\_\_\_ or  $W_s$  30.1  $V_t$  20000  $V_i$  2

Laboratory Name: ATI  
 Case No 6855 / SAS 2726E

Sample Number  
EJ-106

Organics Analysis Data Sheet  
 (Page 4)

Tentatively Identified Compounds

CAS Number	Compound Name	Fraction	RT or Scan Number	Estimated Concentration (ug/l or ug/kg) <i>Sum 2/2/18</i>
1. <u>—</u>	<u>NONE FOUND</u>	<u>VOA</u>	<u>—</u>	<u>—</u>
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30.				

Laboratory Name: ATI  
Case No 6855 / SAS 2726 E

Sample Number  
EJ 106

Organics Analysis Data Sheet  
(Page 4)

Tentatively Identified Compounds

CAS Number	Compound Name	Fraction	RT or Scan Number	Estimated Concentration (ug/l or ug/kg)
1. —	Unknown	BNA	623	1100J
2.				
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015101408  
**Sample Number**  
**EJ 920**

**Organics Analysis Data Sheet**  
 (Page 1)

Laboratory Name: AT7  
 Lab Sample ID No: EJ 920  
 Sample Matrix: Soil  
 Data Release Authorized By: C. Sites

Case No: 6855/SAS 2726 E  
 QC Report No: \_\_\_\_\_  
 Contract No: 68017014  
 Date Sample Received: 2/19/87

**Volatile Compounds**

Concentration: Low Medium (Circle One)  
 Date Extracted/Prepared: 2/24/87  
 Date Analyzed: 2/24/87  
 Conc/Dil Factor: 300 (8.33) = 1.2 pH 7.02  
 Percent Moisture: (Not Decanted) 16.2

CAS Number		ug/l or ug/Kg (Circle One)
74-87-3	Chloromethane	12U
74-83-9	Bromomethane	
75-01-4	Vinyl Chloride	
75-00-3	Chloroethane	
75-09-2	Methylene Chloride	110 B
67-64-1	Acetone	23 B
75-15-0	Carbon Disulfide	6U
75-35-4	1, 1-Dichloroethene	
75-34-3	1, 1-Dichloroethane	
156-60-5	Trans-1, 2-Dichloroethene	
67-66-3	Chloroform	
107-06-2	1, 2-Dichloroethane	
78-93-3	2-Butanone	12U
71-55-6	1, 1, 1-Trichloroethane	6U
56-23-5	Carbon Tetrachloride	
108-05-4	Vinyl Acetate	12U
75-27-4	Bromodichloromethane	6U

CAS Number		ug/l or ug/Kg (Circle One)
78-87-5	1, 2-Dichloropropane	6U
10061-02-6	Trans-1, 3-Dichloropropene	
79-01-6	Trichloroethene	
124-48-1	Dibromochloromethane	
79-00-5	1, 1, 2-Trichloroethane	
71-43-2	Benzene	45 J
10061-01-5	cis-1, 3-Dichloropropene	6U
110-75-8	2-Chloroethylvinylether	12U
75-25-2	Bromoform	6U
108-10-1	4-Methyl-2-Pentanone	12U
591-78-6	2-Hexanone	
127-18-4	Tetrachloroethene	6U
79-34-5	1, 1, 2, 2-Tetrachloroethane	
108-88-3	Toluene	5 JB
108-90-7	Chlorobenzene	6U
100-41-4	Ethylbenzene	3 JB
100-42-5	Styrene	6U
	Total Xylenes	11 B

**Data Reporting Qualifiers**

For reporting results to EPA, the following results qualifiers are used. Additional flags or footnotes explaining results are encouraged. However, the definition of each flag must be explicit.

- Value** If the result is a value greater than or equal to the detection limit, report the value
- U** Indicates compound was analyzed for but not detected. Report the minimum detection limit for the sample with the U (e.g., 10U) based on necessary concentration/dilution action. (This is not necessarily the instrument detection limit.) The footnote should read: U-Compound was analyzed for but not detected. The number is the minimum attainable detection limit for the sample
- J** Indicates an estimated value. This flag is used either when estimating a concentration for tentatively identified compounds where a 1:1 response is assumed or when the mass spectral data indicated the presence of a compound that meets the identification criteria but the result is less than the specified detection limit but greater than zero (e.g., 10J). If limit of detection is 10 ug/l and a concentration of 3 ug/l is calculated, report as 3J
- C** This flag applies to pesticide parameters where the identification has been confirmed by GC/MS. Single component pesticides ≥ 10 ng/l in the final extract should be confirmed by GC/MS
- B** This flag is used when the analyte is found in the blank as well as a sample. It indicates possible/probable blank contamination and warns the data user to take appropriate action
- Other** Other specific flags and footnotes may be required to properly define the results. If used, they must be fully described and such description attached to the data summary report

Laboratory Name AA7  
 Case No: 6855/895 2726 E

Sample Number  
EJ 920

Organics Analysis Data Sheet  
 (Page 2)

Semivolatile Compounds

Concentration: Low Medium (Circle One)  
 Date Extracted/Prepared 2/19/87  
 Date Analyzed: 2/23/87  
 Conc./Dil Factor: 3.26ml / 326g (8.38) x 10<sup>3</sup> = x 200  
 Percent Moisture (Decanted) \_\_\_\_\_

GPC Cleanup  Yes  No  
 Separatory Funnel Extraction  Yes  
 Continuous Liquid-Liquid Extraction  Yes

CAS Number		ug/l or ug/Kg (Circle One)
106-95-2	Phenol	2000u
111-44-4	bis: 2-Chloroethylether	
95-57-8	2-Chlorophenol	
541-73-1	1,3-Dichlorobenzene	
106-46-7	1,4-Dichlorobenzene	
100-51-6	Benzyl Alcohol	
95-50-1	1,2-Dichlorobenzene	
95-48-7	2-Methylnaphenol	
39536-32-9	bis: 2-chloroisopropylether	
106-44-5	4-Methylnaphenol	
621-64-7	N-Nitroso-Di-n-Propylamine	
67-72-1	Hexachloroethane	
98-95-3	Nitrobenzene	
78-59-1	Isophorone	830J
88-75-5	2-Nitrophenol	2000u
105-67-9	2,4-Dimethylnaphenol	
65-85-0	Benzoic Acid	10,000u
111-91-1	bis: 2-ChloroethoxyMethane	2000u
120-83-2	2,4-Dichlorophenol	
120-82-1	1,2,4-Trichlorobenzene	
91-20-3	Naphthalene	
106-47-8	4-Chloroaniline	
87-68-3	Hexachlorobutadiene	
59-50-7	4-Chloro-3-Methylnaphenol	
91-57-6	2-Methylnaphthalene	
77-47-4	Hexachlorocyclopentadiene	
88-06-2	2,4,6-Trichlorophenol	
95-95-4	2,4,5-Trichlorophenol	10,000u
91-58-7	2-Chloronaphthalene	2,000u
88-74-4	2-Nitroaniline	10,000u
131-11-3	Dimethyl Phthalate	2,000u
208-96-8	Acenaphthylene	
99-09-2	3-Nitroaniline	10,000u

CAS Number		ug/l or ug/Kg (Circle One)
83-32-9	Acenaphthene	2000u
51-28-5	2,4-Dinitrophenol	10,000u
100-02-7	4-Nitrophenol	
132-64-9	Dibenzofuran	2000u
121-14-2	2,4-Dinitrotoluene	
606-20-2	2,6-Dinitrotoluene	
84-66-2	Dibenzophthalate	
7005-72-3	4-Chlorophenyl-pnenvietner	
86-73-7	Fluorene	
100-01-6	4-Nitroaniline	10,000u
534-52-1	4,6-Dinitro-2-Methylnaphenol	
86-30-6	N-Nitrosodiphenylamine (1)	2,000u
101-55-3	4-Bromocheny-pnenvietner	
118-74-1	Hexachlorobenzene	
87-86-5	Pentachlorophenol	10,000u
85-01-8	Phenanthrene	800J
120-12-7	Anthracene	2000u
84-74-2	Di-n-Butylphthalate	1300J
206-44-0	Fluoranthene	770J
129-00-0	Pyrene	540J
85-68-7	Butylbenzophthalate	2000u
91-94-1	3,3-Dichlorobenzidine	4000u
56-55-3	Benzofluoranthene	1450J
117-81-7	bis: 2-Ethylhexylphthalate	2000
218-01-9	Chrysene	540J
117-84-0	Di-n-Octyl Phthalate	2000u
205-99-2	Benzofluoranthene	1350J
207-08-9	Benzofluoranthene	2000u
50-32-8	Benzofluoranthene	1560J
193-39-5	Indeno[1,2,3-cd]Pyrene	2000u
53-70-3	Dibenzofluoranthene	
191-24-2	Benzofluoranthene	

(1)-Cannot be separated from diphenylamine

Laboratory Name ATI  
 Case No 6855 / SAS 2726E

Sample Number  
1:10  
EJ 920

Organics Analysis Data Sheet  
 (Page 3)

Pesticide/PCBs

Concentration: Low Medium (Circle One)  
 Date Extracted/Prepared: 19 FEB 87  
 Date Analyzed: 4 MARCH 87  
 Conc/Dil Factor: 0.1  
 Percent Moisture (decanted) 16.2

GPC Cleanup  Yes  No  
 Separatory Funnel Extraction  Yes  
 Continuous Liquid - Liquid Extraction  Yes

CAS Number		ug/l or ug/Kg (Circle One)
319-84-6	Alpha-BHC	94m
319-85-7	Beta-BHC	
319-86-8	Delta-BHC	
58-89-9	Gamma-BHC (Lindane)	
76-44-8	Heptachlor	
309-00-2	Aldrin	
1024-57-3	Heptachlor Epoxide	
959-98-8	Endosulfan I	
60-57-1	Dieldrin	190m
72-55-9	4, 4'-DDE	
72-20-8	Endrin	
33213-65-9	Endosulfan II	
72-54-8	4, 4'-DDD	
1031-07-8	Endosulfan Sulfate	
50-29-3	4, 4'-DDT	
72-43-5	Metoxychlor	940m
53494-70-5	Endrin Ketone	190m
57-74-9	Chlordane	940m
8001-35-2	Toxaphene	1900m
12674-11-2	Aroclor-1016	940m
11104-28-2	Aroclor-1221	
11141-16-5	Aroclor-1232	
53469-21-9	Aroclor-1242	
12672-29-6	Aroclor-1248	
11097-69-1	Aroclor-1254	1900m
11096-82-5	Aroclor-1260	

$V_i$  = Volume of extract injected (ul)

$V_s$  = Volume of water extracted (ml)

$W_s$  = Weight of sample extracted (g)

$V_t$  = Volume of total extract (ul)

$V_s$  \_\_\_\_\_ or  $W_s$  30.6  $V_i$  20000  $V_t$  2

Laboratory Name: ATI  
 Case No 6855 / SAS 2726 E

Sample Number  
EJ-920

Organics Analysis Data Sheet  
 (Page 4)

Tentatively Identified Compounds

CAS Number	Compound Name	Fraction	RT or Scan Number	Estimated Concentration (ug/l or ug/kg)
1. —	NONE FOUND	VOA	—	—
2.				
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Laboratory Name: ATT  
 Case No 6855/875 2726 E

Sample Number  
EJ920

Organics Analysis Data Sheet  
 (Page 4)

Tentatively Identified Compounds

CAS Number	Compound Name	Fraction	RT or Scan Number	Estimated Concentration (ug/l or ug/kg)
1. —	Unknown Xylene	BNA	418	1000 J
2. —	Unknown Xylene	"	459	800 J
3. —	Unknown branched Phenol		1233	1200 J
4. —	Unknown		1548	2600 J
5. —	Unknown		1660	1400 J
6. —	Unknown		1768	13000 J
7. —	Unknown		2026	400 J
8. —	Unknown		2199	3200 J
9. —	Unknown		2268	3000 J
10. —	Unknown		2387	2600 J
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87FPO1559 X  
 Sample Number  
 EJ-104

Organics Analysis Data Sheet  
 (Page 1)

Laboratory Name: ATI  
 Lab Sample ID No: EJ-104  
 Sample Matrix: SOIL  
 Data Release Authorized By: C. Sites

Case No: 6855/SAS 2726E  
 QC Report No: —  
 Contract No: 6801-7014  
 Date Sample Received: 2/19/87

Volatile Compounds

Concentration: Low Medium (Circle One)  
 Date Extracted/Prepared: 2-24-87  
 Date Analyzed: 2-24-87 ✓  
 Conc/Dil Factor: 50 (870) = 1.1 pH 7.45  
 Percent Moisture: (Not Decanted) 13.0

CAS Number		ug/l or ug/Kg (Circle One)
74-87-3	Chloromethane	11U
74-83-9	Bromomethane	
75-01-4	Vinyl Chloride	
75-00-3	Chloroethane	
75-09-2	Methylene Chloride	9B ✓
67-64-1	Acetone	17B ✓
75-15-0	Carbon Disulfide	5U
75-35-4	1, 1-Dichloroethene	
75-34-3	1, 1-Dichloroethane	
156-60-5	Trans-1, 2-Dichloroethene	
67-66-3	Chloroform	
107-06-2	1, 2-Dichloroethane	
78-93-3	2-Butanone	11U
71-55-6	1, 1, 1-Trichloroethane	5U
56-23-5	Carbon Tetrachloride	
108-05-4	Vinyl Acetate	11U
75-27-4	Bromodichloromethane	5U

CAS Number		ug/l or ug/Kg (Circle One)
78-87-5	1, 2-Dichloropropane	5U
10061-02-6	Trans-1, 3-Dichloropropene	
79-01-6	Trichloroethene	
124-48-1	Dibromochloromethane	
79-00-5	1, 1, 2-Trichloroethane	
71-43-2	Benzene	41H ✓
10061-01-5	cis-1, 3-Dichloropropene	5U
110-75-8	2-Chloroethylvinylether	11U
75-25-2	Bromoform	5U
108-10-1	4-Methyl-2-Pentanone	11U
591-78-6	2-Hexanone	
127-18-4	Tetrachloroethene	5U
79-34-5	1, 1, 2, 2-Tetrachloroethane	
108-88-3	Toluene	176B ✓
108-90-7	Chlorobenzene	5U
100-41-4	Ethylbenzene	188B ✓
100-42-5	Styrene	5U
	Total Xylenes	23B

Data Reporting Qualifiers

For reporting results to EPA, the following results qualifiers are used. Additional flags or footnotes explaining results are encouraged. However, the definition of each flag must be explicit.

- Value** If the result is a value greater than or equal to the detection limit, report the value
- U** Indicates compound was analyzed for but not detected. Report the minimum detection limit for the sample with the U (e.g., 10U) based on necessary concentration/dilution action. (This is not necessarily the instrument detection limit.) The footnote should read U-Compound was analyzed for but not detected. The number is the minimum attainable detection limit for the sample
- J** Indicates an estimated value. This flag is used either when estimating a concentration for tentatively identified compounds where a 1:1 response is assumed or when the mass spectral data indicated the presence of a compound that meets the identification criteria but the result is less than the specified detection limit but greater than zero (e.g., 10J). If limit of detection is 10 µg/l and a concentration of 3 µg/l is calculated, report as 3J
- C** This flag applies to pesticide parameters where the identification has been confirmed by GC/MS. Single component pesticides ≥ 10 ng/ul in the final extract should be confirmed by GC/MS
- B** This flag is used when the analyte is found in the blank as well as a sample. It indicates possible/probable blank contamination and warns the data user to take appropriate action
- Other** Other specific flags and footnotes may be required to properly define the results. If used, they must be fully described and such description attached to the data summary report

Laboratory Name ATI  
 Case No. 6855/SAS 2726 E

Sample Number  
EJ104

Organics Analysis Data Sheet  
 (Page 2)

Semivolatile Compounds

Concentration: Low Medium (Circle One)  
 Date Extracted/Prepared 2/19/87 ✓  
 Date Analyzed: 2/22/87  
 Conc./Dil Factor: 30.79(870) × 10<sup>3</sup> = ×37  
 Percent Moisture (Decanted) \_\_\_\_\_

GPC Cleanup  Yes  No  
 Separatory Funnel Extraction  Yes  
 Continuous Liquid-Liquid Extraction  Yes

CAS Number		ug/l or ug/Kg (Circle One)
108-95-2	Phenol	370U
111-44-4	bis(2-Chloroethyl)Ether	
95-57-8	2-Chlorophenol	
541-73-1	1,3-Dichlorobenzene	
106-46-7	1,4-Dichlorobenzene	
100-51-6	Benzyl Alcohol	
95-50-1	1,2-Dichlorobenzene	
95-48-7	2-Methoxyphenol	
39638-32-9	bis(2-chloroisopropyl)Ether	
106-44-5	4-Methoxyphenol	
621-64-7	N-Nitroso-Di-n-Propylamine	
67-72-1	Hexachlorocyclohexane	
98-95-3	Nitrobenzene	
78-59-1	Isophorone	
88-75-5	2-Nitrophenol	
105-67-9	2,4-Dimethylphenol	↓
65-85-0	Benzoic Acid	1800U
111-91-1	bis(2-Chloroethoxy)Methane	370U
120-83-2	2,4-Dichlorophenol	
120-82-1	1,2,4-Trichlorobenzene	
91-20-3	Naphthalene	
106-47-8	4-Chloroaniline	
87-68-3	Hexachlorobutadiene	
59-50-7	4-Chloro-3-Methoxyphenol	↓
91-57-6	2-Methylnaphthalene	62J ✓
77-47-4	Hexachlorocyclopentadiene	370U
88-06-2	2,4,6-Trichlorophenol	↓
95-95-4	2,4,5-Trichlorophenol	1800U
91-58-7	2-Chloronaphthalene	370U
88-74-4	2-Nitroaniline	1800U
131-11-3	Dimethyl Phthalate	370U
208-96-8	Acenaphthylene	↓
99-09-2	3-Nitroaniline	1800U

CAS Number		ug/l or ug/Kg (Circle One)
83-32-9	Acenaphthene	370U
51-28-5	2,4-Dinitrophenol	1800U
100-02-7	4-Nitrophenol	✓ 72J
132-64-9	Dibenzofuran	370U
121-14-2	2,4-Dinitrotoluene	
606-20-2	2,6-Dinitrotoluene	
84-66-2	Dienylphthalate	
7005-72-3	4-Chlorophenylphenylether	
86-73-7	Fluorene	↓
100-01-6	4-Nitroaniline	1800U
534-52-1	4,6-Dinitro-2-Methoxyphenol	370U
86-30-6	N-Nitrosodiphenylamine (1)	140J
101-55-3	4-Bromophenylphenylether	370U
118-74-1	Hexachlorobenzene	↓
87-86-5	Pentachlorophenol	✓ 130J
85-01-8	Phenanthrene	✓ 180J
120-12-7	Anthracene	370U
84-74-2	Di-n-Butylphthalate	✓ 1300U
206-44-0	Fluoranthene	✓ 140J
129-00-0	Pyrene	120J
85-68-7	Butylbenzylphthalate	370U
91-94-1	3,3-Dichlorobenzidine	740U
56-55-3	Benzofluoranthene	110J
117-81-7	bis(2-Ethylhexyl)Phthalate	✓ 160J
218-01-9	Chrysene	✓ 120J
117-84-0	Di-n-Octyl Phthalate	370U
205-99-2	Benzofluoranthene	✓ 170J
207-08-9	Benzofluoranthene	370U
50-32-8	Benzofluoranthene	100J
193-39-5	Indeno[1,2,3-cd]Pyrene	370U
53-70-3	Dibenzofluoranthene	
191-24-2	Benzofluoranthene	↓

(1)-Cannot be separated from diphenylamine

Laboratory Name ATI  
 Case No 6855/SAS 2726 E

Sample Number  
1:5 DIL  
EJ 104

Organics Analysis Data Sheet  
 (Page 3)

Pesticide/PCBs

Concentration: Low Medium (Circle One)  
 Date Extracted/Prepared: 19 FEB 87  
 Date Analyzed: 5 MARCH 87  
 Conc/Dil Factor: 0.2  
 Percent Moisture (decanted) 13.0

GPC Cleanup  Yes  No  
 Separatory Funnel Extraction  Yes  
 Continuous Liquid - Liquid Extraction  Yes

CAS Number		ug/l or ug/Kg (Circle One)
319-84-6	Alpha-BHC	45M
319-85-7	Beta-BHC	
319-86-8	Delta-BHC	
58-89-9	Gamma-BHC (Lindane)	
76-44-8	Heptachlor	
309-00-2	Aldrin	
1024-57-3	Heptachlor Epoxide	
959-98-8	Endosulfan I	↓
60-57-1	Dieldrin	90M
72-55-9	4, 4'-DDE	
72-20-8	Endrin	
33213-65-9	Endosulfan II	
72-54-8	4, 4'-DDD	
1031-07-8	Endosulfan Sulfate	
50-29-3	4, 4'-ODT	↓
72-43-5	Methoxychlor	450M
53494-70-5	Endrin Ketone	90M
57-74-9	Chlordane	450M
8001-35-2	Toxaphene	900M
12674-11-2	Aroclor-1016	450M
11104-28-2	Aroclor-1221	
11141-16-5	Aroclor-1232	
53469-21-9	Aroclor-1242	
12672-29-6	Aroclor-1248	↓
11097-69-1	Aroclor-1254	900M
11096-82-5	Aroclor-1260	↓

$V_i$  = Volume of extract injected (ul)  
 $V_s$  = Volume of water extracted (ml)  
 $W_s$  = Weight of sample extracted (g)  
 $V_t$  = Volume of total extract (ul)

$V_s$  \_\_\_\_\_ or  $W_s$  30.7  $V_t$  20000  $V_i$  2

Laboratory Name: ATI  
Case No 6855/SAS 2726E

Sample Number  
EJ-104

Organics Analysis Data Sheet  
(Page 4)

Tentatively Identified Compounds

CAS Number	Compound Name	Fraction	RT or Scan Number	Estimated Concentration (ug/l or ug/kg)
1. —	NONE FOUND	VOA	—	—
2.				
3.				
4.				
5.				
6.				
7.				
8.				
9.				
10.				
11.				
12.				
13.				
14.				
15.				
16.				
17.				
18.				
19.				
20.				
21.				
22.				
23.				
24.				
25.				
26.				
27.				
28.				
29.				
30.				

Laboratory Name: ATT  
 Case No 6855/SAS 2726 E

Sample Number  
EJ104

Organics Analysis Data Sheet  
 (Page 4)

Tentatively Identified Compounds

CAS Number	Compound Name	Fraction	RT or Scan Number	Estimated Concentration (ug/l or ug/kg)
1. —	unknown chlorinated hydrocarbon	BNA	516	520 J
2. —	unknown aldehyde		587	330 J
3. —	unknown		835	780 J
4. —	unknown phthalate ester		953	480 J
5. —	unknown hydrocarbon		960	560 J
6. —	unk. branched blend		1233	700 J
7. —	unk. branched hydrocarbon		1406	330 J
8. —	unk. branched hydrocarbon		1747	590 J
9. —	unk. branched blend		1914	810 J
10. —	unknown hydrocarbon		2050	370 J
11. —	unk. C <sub>2</sub> =16 hydrocarbon		2113	2200 J
12. —	unk. hydrocarbon		2134	1300 J
13. —	unk. hydrocarbon		2173	1200 J
14. —	unk. hydrocarbon C <sub>2</sub> =35		2242	13,000 J
15. —	unknown		2259	8100 J
16. —	unk. hydrocarbon		2334	1300 J
17. —	unknown		2350	1800 J
18. —	unknown		2419	5500 J
19. —	unknown		2503	3300 J
20. —	unknown		2547	2900 J
21.				
22.				
23.				
24.				
25.				
26.				
27.				
28.				
29.				
30.				

INTRODUCTION TO DATA TABLES

A SUMMARY OF THE ANALYTICAL RESULTS FOR SAMPLES WHICH WERE TAKEN DURING FIELD ACTIVITIES CAN BE FOUND IN THE FOLLOWING TABLES. ONLY DETECTABLE CONCENTRATIONS ARE REPORTED, HOWEVER, IF THE COMPOUND HAS A FOOTNOTE FOLLOWING THE VALUE, CONSULT THE DEFINITION OF THE FOOTNOTE PROVIDED BELOW. ADDITIONAL QA/QC INFORMATION IS PROVIDED IN THE ATTACHED DATA SHEETS.

I) REPORTING UNITS

A) ORGANICS

- 1) Water Samples - ug/l or ppb (parts per billion)
- 2) Soils or Sediments - ug/kg or ppb (parts per billion)

B) METALS

- 1) Water Samples - ug/l or ppb
- 2) Soils or sediments - mg/kg or ppm

II) DEFINITION OF FOOTNOTES TO ANALYTICAL DATA

A) ORGANICS

Footnote	Definition	Interpretation
UJ	Detection Limit (D.L.) is estimated because of a Quality Control (QC) protocol. D.L. is possibly above or below Contract Required Detection Limit (CRDL).	Compound was not detected
UB	Compound found in laboratory blank. No Value above CRDL.	Compound was not detected
UJB	Compound found in laboratory blank, but not detected in sample. CRDL is estimated because of a QC protocol.	Compound was not detected
B	Compound found in blank. Two interpretations are possible: a) If sample value is equivalent to D.L. to 5x blank concentration b) If sample value is greater than 5x the blank concentration	Compound value is semi-quantitative. Compound value is quantitative
JB	Compound found in blank, value is estimated because of QC protocol.	Compound value is semi-quantitative
R	Do Not Use Value. Major Violation of QC Protocol	Compound value is not usable.
C	Value adjusted for blank (an unacceptable procedure)	Compound value is semi-quantitative
J	Value is above CRDL and is an estimated value because of a QC protocol	Compound value is semi-quantitative
Q	No Analytical Result	Compound was not detected
N	Presumptive evidence for the presence of a compound as used for a Tentatively Identified Compound (TIC)	Compound value is semi-quantitative

B) METALS

FOOTNOTE	DEFINITION	INTERPRETATION
E	Estimated or not reported due to interference. See laboratory narrative.	Compound or element was not detected or value is semi-quantitative
s	Analysis by Method of Standard Additions (Look for a "+" Footnote)	Value is quantative
R	Spike recoveries outside QC protocols which indicates a possible matrix problem. Data may be biased high or low. See spike results and laboratory narrative.	Value may be quantitative or semiquantitative
*	Duplicate value outside QC protocols which indicates a possible matrix problem	Value is semiquantitative
+	Correlation coefficient for standard additions is less than 0.995. See review and laboratory narrative.	Data value is biased
→ [ ]	Value is real, but is above instrument D.L. and below CRDL	Value may be quantitative or semiquantitative
UJ	D.L. is estimated because of a QC protocol. D.L. is possibly above or below CRDL.	Compound or element was not detected
J	Value is above CRDL and is an estimated value because of a QC protocol.	Value is semiquantitative

CASE: CR55  
5852726F

VOL  
 ↓  
 Y

	1	2	3	4	5	6	7	8	9	10
IC	EJ	EJ	EJ	EJ						
SAMPLE DTC	104	105	106	920						
COMPOUND										
chloroethane										
bromoethane										
vinyl chloride										
chloroethane										
methylene chloride	9 B	13,000 B	7 B	10 B						
acetone	77 B	5600 B	1 B	23 B						
carbon disulfide										
1,1-dichloroethane		4100 T								
1,1-dichloroethane		8500 T								
trans-1,2-dichloroethane		3000 T								
chloroform										
1,2-dichloroethane										
2-butanone		5100 B								
1,1,1-trichloroethane			17							
carbon tetrachloride										
vinyl acetate										
bromodichloroethane										
1,1,2,2-tetrachloroethane										
1,2-dichloropropane										
trans-1,2-dichloropropene										
trichloroethane		5900 T								
1,1,2-trichloroethane										
benzene		(HL) 7200 T	3 T	4 T						
cis-1,2-dichloropropene										
2-chloroethylvinylether			3 T B							
propofol										
2-hexanone										
4-methyl-2-pentanone										
tetrachloroethane		4600 T								
toluene	16 B	15000 B	4 T B	5 T B						
chlorobenzene	8 B	13000 B	3 T B	3 T B						
ethylbenzene										
styrene										
total xylenes	23 B	110000 B	12 B	11 B						
N-nitrosodimethylamine										
phenol										
aniline										
di(2-chloroethyl)ether										
2-chlorophenol										
1,3-dichlorobenzene										
1,4-dichlorobenzene										
benzyl alcohol										
1,2-dichlorobenzene		(11000)								
2-methylphenol										
di(2-chloroisopropyl)ether										
4-methylphenol										
N-nitroso-di-n-propylamine										
hexachloroethane										
nitrobenzene										
isoprene					8100					
2-nitrophenol										
2,4-dimethylphenol										
benzoic acid										
bis(2-chloroethoxy)ethane										
2,4-dichlorophenol										
1,2,4-trichlorobenzene										
naphthalene		(20000)								
4-chloroaniline										
hexachlorobutadiene										
4-chloro-3-methylphenol										
2-methylnaphthalene		6200 (10000)								
hexachlorocyclopentadiene										
2,4,6-trichlorophenol										
2,4,5-trichlorophenol										
2-chloronaphthalene										
2-nitroaniline										
dimethyl phthalate										
acenaphthylene										
3-nitroaniline		(7000)								
acenaphthene										
2,4-dinitrophenol										
4-nitrophenol										
dibenzofuran										
2,4-dinitrotoluene										
2,6-dinitrotoluene										
diethylphthalate										
4-chlorophenyl-phenylether										
flourene										
4-nitroaniline										
4,6-dinitro-2-methylphenol										
N-nitrosodiphenylamine		1400	(20000)							
4-bromophenyl-phenylether										
hexachlorobenzene										

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= Use for HRS  
 = Tini quatation  
 = unmarked means  
 an artifact  
 etc

B = found in  
 blank  
 J = estimate

R = do not use  
 calibration  
 problem

CASE: 6055  
SAS 2726E

1 2 3 4 5 6 7 8 9 10

COMPOUND	1	2	3	4	5	6	7	8	9	10
pentachlorophenol	130J	55000								
phenanthrene	180J	7500J		800J						
anthracene										
di-n-butylphthalate	1900		1100J							
fluoranthene	140J		240J	990J						
benzidine				540J						
pyrene	120J		250J							
butylbenzylphthalate										
3,3'-dichlorobenzidine										
benzo(a)anthracene	110J			450J						
bis(2-ethylhexyl)phthalate		100,000	14000							
chrysene	120J			540J						
di-n-octyl phthalate										
benzo(b)fluoranthene	170J			550J						
benzo(a)pyrene	100J			560J						
indeno(1,2,3-cd)pyrene										
dibenzo(a,h)anthracene										
benzo(g,h,i)perylene										
alpha-BHC										
beta-BHC										
delta-BHC										
gamma-BHC(lindane)										
heptachlor			BDL							
aldrin										
heptachlor epoxide										
endosulfan I										
dieldrin										
4,4'-DDE										
endrin										
endosulfan II										
4,4'-DDD										
endrin aldehyde										
endosulfan sulfate										
4,4'-DDT										
methoxychlor										
endrin ketone										
chlorodane										
carbazene										
Aroclor-1016										
Aroclor-1221										
Aroclor-1232										
Aroclor-1242										
Aroclor-1248										
Aroclor-1254										
Aroclor-1260										
ELMENT										
aluminum										
antimony										
arsenic										
barium										
beryllium										
caesium										
calcium										
chromium										
cobalt										
copper										
iron										
lead										
magnesium										
manganese										
mercury										
nickel										
potassium										
selenium										
silver										
sodium										
thallium										
tin										
vanadium										
zinc										
cyanide CHECK IF ANALYZED ( )										
TENTATIVELY IDENTIFIED ORGANICS										

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BDL  
Below  
detection  
Limit

O = use for HRS  
□ = semi  
quantative  
= unmarked  
means an  
artifact  
TC

5F3778



Analytical **Technologies, Inc.**

Corporate Offices: 5550 Morehouse Drive San Diego, CA 92121 (619) 458-9141

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Analytical Technologies, Inc.  
5550 Morehouse Drive  
San Diego, CA 92121

March 5, 1987

<sup>6855</sup>  
Case Narrative for EPA Case ~~8655~~/SAS 2726E  
Contract No: 6801-7014

Soil Samples: EJ-104, EJ-105, EJ-106, EJ-920

This case was received at Analytical Technologies, Inc.  
on 2/19/87. Comments regarding the analytical fractions  
are contained in the following attachments.

Sincerely,

*Robert V. Woods*  
Robert V. Woods

RVW/sm  
Attachments

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MAR 6 1987

U.S. EPA, CENTRAL REGIONAL OFFICE  
536 S. CLARK STREET  
CHICAGO, ILLINOIS 60605

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Attachment 1

General Notes

1. GC/MS files that have a M1 in the name denote analyses were done using GC/MS system #1. 'M2' and 'M3' are used to identify files acquired using GC/MS systems #2 and #3 respectively.
2. The GC/MS dual display of the HSL 'hit' and the library standard spectra with the difference between the two is not a comparison with the NBS Library but rather a library created using spectra obtained on the specific GC/MS instrument.
3. GC/MS 'TIC' compounds are calculated using the appropriate internal standard's peak height (measured by caliper and converted to RIC units). The 'TIC' compound's peak height used in calculations is the RIC enhanced value displayed on the forward library search printout. The calculations written on the p/o's are for ng/ul of the TIC compound in the extract. The formula written is typically as follows:

$$\frac{(\text{RIC height of TIC}) * (\text{ISTD amount})}{(\text{RIC peak height Istd}) * (\text{response factor}=1)} = \text{ng/ul TIC}$$

4. On the primary GC column, 'NC' refers to a compound not confirmed on the confirmation column.
5. For the the confirmation column, 'NPR' refers to a compound not found during the primary run.

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CASE 6855/SAS 2726E  
COMMENTS ON VOA ANALYSES

- 1) Soil sample EJ-105 was analyzed as a medium level soil in accordance with CLP medium level soil extraction.
  - a. 10 uL of the extract was purged with 5 mL water.
  - b. Surrogate values listed on the quantitation report must be multiplied by a factor of 10. All surrogate recoveries met QA/QC limits.
  - c. Spike values listed on the quantitation report also must be multiplied by a factor of 10. All spike recoveries were low. Benzene recoveries did not meet QA/QC criteria for both MS and MSD. Recoveries for tri-chloroethene and toluene were low in the MS sample. The low recoveries may be due to the required sample dilution, which lowered the spike concentration levels close to the instrument detection limits. For some of the compounds, the amount spiked was much lower than the sample concentrations which may have contributed to the low recoveries.
  - d. The medium level methanol blank contained chloroform, ethyl benzene and xylene at or above the CRDL. The methanol used for the sample extractions was the least contaminated available to us at that time. We have notified the manufacturer and they are aware of the problem. They are sending us a new lot of methanol which is supposed to be less contaminated, but it was not possible to wait for that methanol without exceeding the VOA holding time.

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COMMENTS ON BNA ANALYSES

- 1) One of the BNA 5-point calibration curves included standards analyzed on two different days. There were four (4) standards analyzed on February 20, 1987, and 50 Ng Standard was analyzed on February 21, 1987. This 5-point curve is referenced as February 21, 1987.
- 2) The continuing calibration standards analyzed on February 23, 1987, and February 28, 1987, exceeded R.f. criteria for a C.C.C. Compound, N-Nitrosodiphenylamine. We attempted to fix the problem by remaking the standard, changing the injection port liner, and removing the front section of the column. However, this was not successful. This problem affected two (2) samples: EJ106 and EJ920. N-Nitrosodiphenylamine was not found in either sample.
- 3) The Sample EJ920 extract could not be concentrated below 5 mls due to the high concentration of high molecular weight compounds. The BNA chromatograms shows a high concentration of hydrocarbons in the range of C<sub>18</sub> to ≥ C<sub>28</sub>.

Comments on Pesticide Analysis:

On 3/3/87 most components in IND Mix B CONT 1 (2054) exceeded calibration factor criteria because IND Mix B Initial was not allowed to reach room temperature before it was analyzed. There were not any samples ran after this standard, and the 12 hour sequence was restarted.

On 3/3/87 Endosulfan II in IND A Cont 1 (1711) exceeded calibration factor criteria due to a contaminant which coeluted with this compound. There were not any hits for this compound and no samples were affected.

On 3/3/87 Endosulfan Sulfate in IND B Cont 2 (0448) exceeded calibration factor criteria. However, there were not any samples ran after this sample.

On 3/2/87 the retention times of DDT and DBC in EVAL Mix A (2344) were not labeled by the integrator. Thus the retention time shift for Dibutyl Chlorendate is based upon EVAL Mix B (0038).

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21 DAY CLOCK

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ASE: 2855/SA 5 2726 E

DATE OF START: 12 Mar 87

ITE:

ISSUED BY: C. Eddy

% OF SAMPLES: 4

ACCEPTED BY: [Signature]

STOP NOTICE

DATE OF STOP: 12 Mar 87

ISSUED BY:

REASON:

ACCEPTED BY:

- (1) Friday required field training
- (2) Restart review 16 Mar 87;  
Including training.
- (3) Stopped on 16
- (4) Restarted and completed on 24 Mar 87
- (5) Stopped for typos.

Delays caused by mandatory training  
travel and sickness.

Review time included data review training  
completed w/ 14 days [Signature]  
[Signature]

DATE OF RESTART

ISSUED BY

ACCEPTED BY

DATE OF COMPLETION

ACCEPTED BY

# SCM SUPPOCATE PERCENT RECOVERY SUMMARY

Case No. 68558952726E Contract Laboratory AT7 Contract No. 68017014

Low  Medium

SMD TRAFFIC NO.	VOLATILE				SEMI-VOLATILE				PESTICIDE	
	TOLUENE-08 (99-117)	BFB (74-121)	1,2 DICHLORO-ETHANE-04 (70-121)	NITRO-BENZENE-05 (23-120)	2-FLUORO-BIPHENYL (30-119)	TERPENEYL-DIA (18-127)	PHENOL-05 (24-119)	2-FLUORO-PHENOL (25-121)	2,4,6 TRIBROMO-PHENOL (18-122)	DIBUTYL-CHLOROPHOSPHATE (20-150)
EJ104	111	82	114	87	86	83	74	95	68	89
EJ105	N/A			83	81	100	63	110	91	73
EJ106	100	99	105	39	98	88	*23	47	65	93
EJ120	107	91	111	85	98	93	40	80	121	109
EJ106MS	112	89	113	60	64	104	48	70	82	100
EJ106MSD	106	93	109	52	68	116	48	82	77	106
R2-1982	N/A			73	78	85	71	81	65	107
R2-2487	101	109	103	N/A						N/A

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\* VALUES ARE OUTSIDE OF CONTRACT REQUIRED QC LIMITS Volatiles: 0 out of 18 ; outside of QC limits  
 \* ADVISORY LIMITS ONLY Semi-Volatiles: 1 out of 42 ; outside of QC limits  
 Pesticides: 0 out of 2 ; outside of QC limits

Comments: N/A = Not APPLICABLE

# SO<sup>2</sup> SINTERCATE RECOVERY SUMMARY

Case No. 6855/SAS 2726E Contract Laboratory ATI Contract No. 6801-7014

SNO TRAFFIC NO.	VOLATILE				SEMI-VOLATILE				PESTICIDE	
	TOLUENE-08 (81-117)	BFB (74-121)	1,2 DICHLORO- ETHANE-04 (70-121)	NITRO- BENZENE-09 (83-120)	2-FLUORO- BIPHENYL (50-118)	TERPENEYL- 014 (18-137)	PHENOL-03 (24-113)	2-FLUORO- PHENOL (28-121)	2,4,6 TRIBROMO PHENOL (10-122)	DIBUTYL- CHLOROPHOSPHATE (20-150)
RB <sub>2</sub> MLCH	99	107	101	N/A						N/A
EJ-105	89	107	94							
EJ-105MS	79*	89	86							
EJ-105MISO	83	100	94							

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Volatiles: out of 12 ; outside of QC limits  
 Semi-Volatiles: out of N/A ; outside of QC limits  
 Pesticides: out of N/A ; outside of QC limits

7/85

\* VALUES ARE OUTSIDE OF CONTRACT REQUIRED QC LIMITS

\*\* ADVISORY LIMITS ONLY

Comments: N/A = Not APPLICABLE

**SOIL MATRIX SPIKE/MATRIX SPIKE DUPLICATE RECOVERY**

Case No. 6855/SAS 2726E Contractor ATI Contract No. 6801-7014

Low Level X Medium Level \_\_\_\_\_

FRACTION	COMPOUND	CONC. SPIKE ADDED (ug/Kg)	SAMPLE RESULT	CONC. MS	% REC	CONC. MSD	% REC	RPD	OC LIMITS
VOA SMO SAMPLE NO. EJ-106	1,1-Dichloroethene	50	BDL	50	100	48	96	4	22 59-172
	Trichloroethene		↓	41	82	43	86	5	24 62-137
	Chlorobenzene		↓	51	102	49	98	4	21 60-133
	Toluene		↓	55	102	51	94	8	21 59-139
	Benzene		↓	51	96	49	92	4	21 66-142
B/N SMO SAMPLE NO. EJ106	1,2,4-Trichlorobenzene	175	BDL	146	83	135	77	8	23 38-107
	Acenaphthene	100	↓	78	78	89	89	13	19 31-137
	2,4-Dinitrotoluene		↓	69	69	75	75	8	47 28-89
	Pyrene		↓	118	116	131	129	11	36 35-142
	N-Nitrosodi-n-Propylamine		↓	52	52	48	48	8	38 41-126
ACID SMO SAMPLE NO. EJ106	1,4-Dichlorobenzene	200	↓	77	38	76	38	7	27 28-104
	Pentachlorophenol		↓	99	50	87	44	13	47 17-109
	Phenol		↓	55	28	63	32	13	35 28-90
	2-Chlorophenol		↓	119	60	126	63	5	50 25-102
	4-Chloro-3-Methylphenol		↓	120	60	128	64	8	33 26-103
PEST SMO SAMPLE NO. EJ	4-Nitrophenol		↓	130	65	103	52	22	50 11-114
	Lindane	27.4	↓	28.1	103	28.8	105	1.9	50 46-127
	Heptachlor	27.4	↓	24.5	108	20.8	113	1.5	31 35-130
	Aldrin	27.4	↓	34.9	128	48.2	176	7.2	43 34-132
	Dieldrin	68.6	↓	70.0	102	85.8	125	2.0	38 31-134
EJ	Endrin	68.6	↓	98.6	142	72.6	135	2.2	45 42-139
	4,4'-DDT	68.6	↓	81.6	119	88.5	129	8.1	50 23-134

\*ASTERISKED VALUES ARE OUTSIDE OC LIMITS.

RPD: VOA: 0 out of 5; outside OC limits  
 B/N: 0 out of 6; outside OC limits  
 ACID: 0 out of 5; outside OC limits  
 PEST: 0 out of 6; outside OC limits

RECOVERY: VOA: 0 out of 10; outside OC limits  
 B/N: 0 out of 12; outside OC limits  
 ACID: 0 out of 10; outside OC limits  
 PEST: 1 out of 12; outside OC limits

Comments: BNA Matrix spike concentration determined by comparison with EPA traceable standard

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# SOIL MATRIX SPIKE / MATRIX SPIKE DUPLICATE RECOVERY

Case No. 6855/SAS 2726E

Contractor ATI

Contract No. 68017014

Low Level \_\_\_\_\_ Medium Level X

FRACTION	COMPOUND	CONC. SPIKE ADDED (ug/Kg)	SAMPLE RESULT	CONC. MS	% REC	CONC. MSD	% REC	RPD	OC LIMITS
VOA SMO SAMPLE NO. EJ-105	1,1-Dichloroethene	50	19	54	70	58	78	11	22
	Trichloroethene		27	52	150*	58	62	21	24
	Chlorobenzene		BVL	32	64	37	74	14	21
	Toluene		67	88	122*	97	60	35*	21
	Benzene		33	31	10*	35	44	2	21
B/N SMO SAMPLE NO. N/A	1,2,4-Trichlorobenzene	N/A							
	Acenaphthene								
	2,4-Dinitrotoluene								
	Pyrene								
	N-Nitrosodi-n-Propylamine								
ACID SMO SAMPLE NO. N/A	1,4-Dichlorobenzene								
	Pentachlorophenol								
	Phenol								
	2-Chlorophenol								
	4-Chloro-3-Methylphenol								
PEST SMO SAMPLE NO. N/A	4-Nitrophenol								
	Lindane								
	Heptachlor								
	Aldrin								
	Diieldrin								
	Endrin								
	4,4'-DDT								

\*ASTERISKED VALUES ARE OUTSIDE OC LIMITS.

RPD: VOA: 1 out of 5 : outside OC limits  
 B/N: 150 out of N/A : outside OC limits  
 ACID: 1 out of 1 : outside OC limits  
 PEST: 1 out of 1 : outside OC limits

RECOVERY: VOA: 4 out of 10 : outside OC limits  
 B/N: 150 out of N/A : outside OC limits  
 ACID: 1 out of 1 : outside OC limits  
 PEST: 1 out of 1 : outside OC limits

Comments: N/A = NOT APPLICABLE

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# METHOD BLANK SUMMARY

Case No. 6855 Region 5 Contractor ATI Contract No. 68017014  
SAS2726E

FILE NO	DATE OF ANALYSIS	FRACTION	MATRIX	CONC. LEVEL	INST. ID	CAS NUMBER	COMPOUND (MSL, TIC OR UNKNOWN)	CONC.	UNITS	CRDL
RB 2/19/87	3 MAR 1987	REST	SOIL	L	HPDRA		NONE			
RB 2/19/87	2/22/87	BWA	SOIL	L	M3	87-74-2	Di-n-Butylphthalate	780	ug/kg	330U
							unknown Diphenyl	200J		
							unknown Silane	230J		
							unknown Silane	230J		
							unknown branched phenol	430J		
							unknown phthalate	130J		
							Methylene Chloride	6.	ug/kg	56U
RB 2/24/87	2/24/87	VQA	H2O	L	M1	67-64-1	Acetone	3J		106U
							Benzoforn	3J		56U
							Toluene	3J		56U
							Ethyl benzene	2J		56U
							Total Xylenes	8		56U
							Methylene Chloride	1000		250U
RB 2 MeOH	2/25/87	VQA	H2O	M	M1	75-09-2	Acetone	150J		500U
							Methoforn	250		250U
							2-Butanone	210J		500U
							Toluene	160J		250U
							Ethyl benzene	300		250U
							Total Xylenes	300		250U

Comments:

87FPO1559

Sample Number

EJ-104

RECEIVED MAR 30 1987  
Organics Analysis Data Sheet  
(Page 1)

Laboratory Name: ATI  
Lab Sample ID No: EJ-104  
Sample Matrix: SOIL  
Data Release Authorized By: C. Sites

Case No: 6855/SAS 2726E  
QC Report No: —  
Contract No: 6801-7014  
Date Sample Received: 2/19/87

## Volatile Compounds

Concentration: Low Medium (Circle One)Date Extracted/Prepared: 2-24-87Date Analyzed: 2-24-87Conc/Dil Factor: 50 (870) = 1.1 pH 7.45 1.15Percent Moisture: (Not Decanted) 13.0

$$\frac{100-13}{100} = 87$$

CAS Number		ug/l or ug/Kg (Circle One)
74-87-3	Chloromethane	11U
74-83-9	Bromomethane	
75-01-4	Vinyl Chloride	
75-00-3	Chloroethane	
75-09-2	Methylene Chloride	9B
67-64-1	Acetone	17B
75-15-0	Carbon Disulfide	5U
75-35-4	1, 1-Dichloroethene	
75-34-3	1, 1-Dichloroethane	
156-60-5	Trans-1, 2-Dichloroethene	
67-66-3	Chloroform	
107-06-2	1, 2-Dichloroethane	
78-93-3	2-Butanone	11UR <sup>2</sup> K
71-55-6	1, 1, 1-Trichloroethane	5U
56-23-5	Carbon Tetrachloride	
108-05-4	Vinyl Acetate	11U
75-27-4	Bromodichloromethane	5U

CAS Number		ug/l or ug/Kg (Circle One)
78-87-5	1, 2-Dichloropropane	5U
10061-02-6	Trans-1, 3-Dichloropropene	
79-01-6	Trichloroethene	
124-48-1	Dibromochloromethane	
79-00-5	1, 1, 2-Trichloroethane	
71-43-2	Benzene	141/11U
10061-01-5	cis-1, 3-Dichloropropene	5U
110-75-8	2-Chloroethylvinylether	11UR <sup>2</sup> K
75-25-2	Bromoform	5U
108-10-1	4-Methyl-2-Pentanone	11U
591-78-6	2-Hexanone	
127-18-4	Tetrachloroethene	5U
79-34-5	1, 1, 2, 2-Tetrachloroethane	
108-88-3	Toluene	116B <sup>1</sup>
108-90-7	Chlorobenzene	5U
100-41-4	Ethylbenzene	188 <sup>1</sup>
100-42-5	Styrene	5U
	Total Xylenes	123B <sup>1</sup>

## Data Reporting Qualifiers

For reporting results to EPA, the following results qualifiers are used.  
Additional flags or footnotes explaining results are encouraged. However, the definition of each flag must be explicit.

- Value** If the result is a value greater than or equal to the detection limit, report the value
- U** Indicates compound was analyzed for but not detected. Report the minimum detection limit for the sample with the U (e.g., 10U) based on necessary concentration/dilution factor. (This is not necessarily the instrument detection limit.) The footnote should read: U-Compound was analyzed for but not detected. The number is the minimum attainable detection limit for the sample
- J** Indicates an estimated value. This flag is used either when estimating a concentration for tentatively identified compounds where a 1:1 response is assumed or when the mass spectral data indicated the presence of a compound that meets the identification criteria but the result is less than the specified detection limit but greater than zero (e.g., 10J) If limit of detection is 10 ug/l and a concentration of 3 ug/l is calculated, report as 3J

- C** This flag applies to pesticide parameters where the identification has been confirmed by GC/MS. Single component pesticides  $\geq 10$  ng/l in the final extract should be confirmed by GC/MS
- B** This flag is used when the analyte is found in the blank as well as a sample. It indicates possible/probable blank contamination and warns the data user to take appropriate action
- Other** Other specific flags and footnotes may be required to properly define the results. If used, they must be fully described and such description attached to the data summary report

Laboratory Name ATI  
 Case No: 6855/SAS 2726 E

RECEIVED MAR 30 1987

Sample Number  
EJ104

Organics Analysis Data Sheet  
 (Page 2)

Semivolatile Compounds

Concentration: Low Medium (Circle One)  
 Date Extracted/Prepared: 2/19/87  
 Date Analyzed: 2/22/87  
 Conc./Dil Factor: 30.7g(-870) x 10<sup>3</sup> = x37  
 Percent Moisture (Decanted): \_\_\_\_\_

GPC Cleanup  Yes  No  
 Separatory Funnel Extraction  Yes  
 Continuous Liquid-Liquid Extraction  Yes

CAS Number	Compound	ug/l or ug/Kg (Circle One)
108-95-2	Phenol	370U
111-44-4	bis: 2-Chloroethoxy Ethane	
95-57-6	2-Chlorophenol	
541-73-1	1,3-Dichlorobenzene	
106-46-7	1,4-Dichlorobenzene	
100-51-6	Benzyl Alcohol	URK
95-50-1	1,2-Dichlorobenzene	
95-48-7	2-Methoxyphenol	
39538-32-9	bis: 2-chloroisopropyl Ether	
106-44-5	4-Methoxyphenol	
621-64-7	N-Nitroso-Di-n-Propylamine	
67-72-1	Hexachloroethane	
98-95-3	Nitrobenzene	
78-59-1	Isopropylone	
86-75-5	2-Nitrophenol	
105-67-9	2,4-Dimethoxyphenol	↓
65-85-0	Benzoic Acid	1800U
111-91-1	bis: 2-Chloroethoxy Methane	370U
120-83-2	2,4-Dichlorophenol	
120-82-1	1,2,4-Trichlorobenzene	
91-20-3	Naphthalene	
106-47-8	4-Chloroaniline	REK
87-68-3	Hexachlorocyclopentadiene	
59-50-7	4-Chloro-3-Methoxyphenol	↓
91-57-6	2-Methoxynaphthalene	162J ✓
77-47-4	Hexachlorocyclopentadiene	370U J26
88-06-2	2,4,6-Trichlorophenol	↓
95-95-4	2,4,5-Trichlorophenol	1800U
91-58-7	2-Chloronaphthalene	370U
88-74-4	2-Nitroaniline	1800U
131-11-3	Dimethyl Phthalate	370U
208-96-8	Acenaphthylene	↓
99-09-2	3-Nitroaniline	1800U REK

CAS Number	Compound	ug/l or ug/Kg (Circle One)
83-32-9	Acenaphthene	370U
51-28-5	2,4-Dinitrophenol	1800U
100-02-7	4-Nitrophenol	172J ✓
132-64-9	Dibenzofuran	1370U
121-14-2	2,4-Dinitrotoluene	
606-20-2	2,6-Dinitrotoluene	
84-66-2	Dibenzofuran	
7005-72-3	4-Chlorophenylphenyl ether	
86-73-7	Fluorene	↓
100-01-6	4-Nitroaniline	1800U
534-52-1	4,6-Dinitro-2-Methoxyphenol	1370U
86-30-6	N-Nitrosodiphenylamine (1)	140J ✓
101-55-3	4-Bromophenylphenyl ether	370U
118-74-1	Hexachlorobenzene	↓
87-86-5	Pentachlorophenol	1130J ✓
85-01-8	Phenanthrene	1180J ✓
120-12-7	Anthracene	370U
84-74-2	Di-n-Butylphthalate	11300J ✓
206-44-0	Fluoranthene	1740J ✓
129-00-0	Pyrene	1120J ✓
85-68-7	Butylbenzophthalate	370U
91-94-1	3,3-Dichlorobenzidine	740U J26
56-55-3	Benz[a]Anthracene	1110J ✓
117-81-7	bis: 2-Ethylhexyl Phthalate	160J ✓
218-01-9	Chrysene	1120J ✓
117-84-0	Di-n-Octyl Phthalate	370U
205-99-2	Benz[a]Fluoranthene	1170J ✓
207-08-9	Benz[a]Fluoranthene	370U
50-32-8	Benz[a]Pyrene	1100J ✓
193-39-5	Indeno[1,2,3-cd]Pyrene	370U
53-70-3	Dibenz[a,h]Anthracene	↓
191-24-2	Benz[a,h]Perylene	↓ J26

(1)-Cannot be separated from diphenylamine

Laboratory Name ATI RECEIVED MAR 30 1987  
 Case No 6855/SAS 2726 E

Sample Number  
1:5 DIL  
EJ 104

Organics Analysis Data Sheet  
 (Page 3)

Pesticide/PCBs

Concentration: Low Medium (Circle One)  
 Date Extracted/Prepared: 19 FEB 87  
 Date Analyzed: 5 MARCH 87  
 Conc/Dil Factor: 0.2  
 Percent Moisture (decanted) 13.0

GPC Cleanup  Yes  No  
 Separatory Funnel Extraction  Yes  
 Continuous Liquid - Liquid Extraction  Yes

CAS Number		ug/l or ug/Kg (Circle One)
319-84-6	Alpha-BHC	45m
319-85-7	Beta-BHC	
319-86-8	Delta-BHC	
58-89-9	Gamma-BHC (Lindane)	
76-44-8	Heptachlor	
309-00-2	Aldrin	
1024-57-3	Heptachlor Epoxide	
959-98-8	Endosulfan I	↓
60-57-1	Dieldrin	90m
72-55-9	4,4'-DDE	
72-20-8	Endrin	
33213-65-9	Endosulfan II	
72-54-8	4,4'-DDD	
1031-07-8	Endosulfan Sulfate	
50-29-3	4,4'-DDT	↓
72-43-5	Methoxychlor	450m
53494-70-5	Endrin Ketone	90m
57-74-9	Chlordane	450m
8001-35-2	Toxaphene	900m
12674-11-2	Aroclor-1016	450m
11104-28-2	Aroclor-1221	
11141-16-5	Aroclor-1232	
53469-21-9	Aroclor-1242	
12672-29-6	Aroclor-1248	↓
11097-69-1	Aroclor-1254	900m
11096-82-5	Aroclor-1260	↓

$V_i$  = Volume of extract injected (ul)

$V_s$  = Volume of water extracted (ml)

$W_s$  = Weight of sample extracted (g)

$V_t$  = Volume of total extract (ul)

$V_s$  \_\_\_\_\_ or  $W_s$  30.7  $V_i$  20000  $V_t$  2

RECEIVED MAR 30 1987

Laboratory Name: ATI  
Case No 6855/SAS 2726E

Sample Number  
EJ-104

Organics Analysis Data Sheet  
(Page 4)

Tentatively Identified Compounds

CAS Number	Compound Name	Fraction	RT or Scan Number	Estimated Concentration (ug/l or ug/kg)
1. —	NONE FOUND	VOA	—	—
2.				
3.				
4.				
5.				
6.				
7.				
8.				
9.				
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Laboratory Name: ATT RECEIVED MAR 30 1987  
 Case No 6855/SAS 2726E

Sample Number  
EJ104

Organics Analysis Data Sheet  
 (Page 4)

Tentatively Identified Compounds

CAS Number	Compound Name	Fraction	RT or Scan Number	Estimated Concentration (ug/l or ug/kg)
1. —	unknown chlorinated hydrocarbon	BNA	576	520 J
2. —	unknown aldehyde		587	330 J
3. —	unknown		835	780 J
4. —	unknown phthalate ester		953	480 J
5. —	unknown hydrocarbon		960	560 J
6. —	unk. branched blend		1233	700 J
7. —	unk. branched hydrocarbon		1406	330 J
8. —	unk. branched hydrocarbon		1747	590 J
9. —	unk. branched blend		1914	810 J
10. —	unknown hydrocarbon		2050	370 J
11. —	unk. C <sub>316</sub> hydrocarbon		2113	2200 J
12. —	unk. hydrocarbon		2134	1300 J
13. —	unk. hydrocarbon		2173	1200 J
14. —	unk. hydrocarbon C <sub>335</sub>		2242	13,000 J
15. —	unknown		2259	8100 J
16. —	unk. hydrocarbon		2334	1300 J
17. —	unknown		2350	1800 J
18. —	unknown		2419	5500 J
19. —	unknown		2503	3300 J
20. —	unknown		2547	2900 J
21. —				
22. —				
23. —				
24. —				
25. —				
26. —				
27. —				
28. —				
29. —				
30. —				

87FPO1560

Sample Number  
EJ-105

Organics Analysis Data Sheet  
RECEIVED MAR 30 1987 (Page 1)

Laboratory Name: ATI  
Lab Sample ID No: EJ-105  
Sample Matrix: SOIL  
Data Release Authorized By: C Sites

Case No: 6855/SAS 2726E  
QC Report No: —  
Contract No: 6801-7014  
Date Sample Received: 2-19-87

Volatile Compounds

Concentration: Low Medium (Circle One)

Date Extracted/Prepared: 2-25-87

Date Analyzed: 2-25-87

Conc/Dil Factor:  $\frac{10 \text{ mL}}{4.0 \text{ g}} \cdot \frac{2 \text{ mL}}{0.01 \text{ mL}} \cdot \frac{1}{50} \cdot \frac{2200}{\text{Dil}}$  8.82

Percent Moisture: (Not Decanted) 43.8

CAS Number	Compound	ug/l or ug/Kg (Circle One)
74-87-3	Chloromethane	22,000U
74-83-9	Bromomethane	
75-01-4	Vinyl Chloride	
75-00-3	Chloroethane	
75-09-2	Methylene Chloride	13,000B
67-64-1	Acetone	5,600JB
75-15-0	Carbon Disulfide	11,000U
75-35-4	1, 1-Dichloroethene	4,100J
75-34-3	1, 1-Dichloroethane	8,800J
156-60-5	Trans-1, 2-Dichloroethene	3,000J
67-66-3	Chloroform	11,000U
107-06-2	1, 2-Dichloroethane	
78-93-3	2-Butanone	5,100JBKJ
71-55-6	1, 1, 1-Trichloroethane	200,000
56-23-5	Carbon Tetrachloride	11,000U
108-05-4	Vinyl Acetate	22,000U
75-27-4	Bromodichloromethane	11,000U

CAS Number	Compound	ug/l or ug/Kg (Circle One)
78-87-5	1, 2-Dichloropropane	14,000U
10061-02-6	Trans-1, 3-Dichloropropene	
79-01-6	Trichloroethene	5,900J
124-48-1	Dibromochloromethane	11,000U
79-00-5	1, 1, 2-Trichloroethane	
71-43-2	Benzene	7,200J
10061-01-5	cis-1, 3-Dichloropropene	11,000U
110-75-8	2-Chloroethylvinylether	22,000U
75-25-2	Bromoform	11,000U
108-10-1	4-Methyl-2-Pentanone	22,000U
591-78-6	2-Hexanone	
127-18-4	Tetrachloroethene	14,600J
79-34-5	1, 1, 2, 2-Tetrachloroethane	11,000U
108-88-3	Toluene	15,000B
108-90-7	Chlorobenzene	11,000U
100-41-4	Ethylbenzene	33,000B
100-42-5	Styrene	11,000U
	Total Xylenes	160,000B

Data Reporting Qualifiers

For reporting results to EPA, the following results qualifiers are used. Additional flags or footnotes explaining results are encouraged. However, the definition of each flag must be explicit.

- Value** If the result is a value greater than or equal to the detection limit, report the value
- U** Indicates compound was analyzed for but not detected. Report the minimum detection limit for the sample with the U (e.g., 10U) based on necessary concentration/dilution action. (This is not necessarily the instrument detection limit.) The footnote should read: U-Compound was analyzed for but not detected. The number is the minimum attainable detection limit for the sample
- J** Indicates an estimated value. This flag is used either when estimating a concentration for tentatively identified compounds where a 1:1 response is assumed or when the mass spectral data indicated the presence of a compound that meets the identification criteria but the result is less than the specified detection limit but greater than zero. (e.g., 10J) If limit of detection is 10 ug/l and a concentration of 3 ug/l is calculated, report as 3J

- C** This flag applies to pesticide parameters where the identification has been confirmed by GC/MS. Single component pesticides ≥ 10 ng/ul in the final extract should be confirmed by GC/MS
- B** This flag is used when the analyte is found in the blank as well as a sample. It indicates possible/probable blank contamination and warns the data user to take appropriate action
- Other** Other specific flags and footnotes may be required to properly define the results. If used, they must be fully described and such description attached to the data summary report

Laboratory Name AT7  
 Case No. 6855/898 2726 E

Sample Number  
EJ 105

Organics Analysis Data Sheet  
 (Page 2)

RECEIVED MAR 30 1987

Semivolatile Compounds

Concentration: Low Medium (Circle One)  
 Date Extracted/Prepared 2/19/87  
 Date Analyzed: 3/3/87  
 Conc./Dil Factor: 5.26ml / 30.505(562) x 10<sup>3</sup> x 610  
 Percent Moisture (Decanted) \_\_\_\_\_

GPC Cleanup  Yes  No  
 Separatory Funnel Extraction  Yes  
 Continuous Liquid-Liquid Extraction  Yes

CAS Number	Compound	ug/l or ug/Kg (Circle One)
108-95-2	Phenol	6100U
111-44-4	bis(2-Chloroethoxy)Ethane	J
95-57-8	2-Chlorophenol	
541-73-1	1,3-Dichlorobenzene	
106-46-7	1,4-Dichlorobenzene	
100-51-6	Benzyl Alcohol	
95-50-1	1,2-Dichlorobenzene	11,000
95-48-7	2-Methoxyphenol	6100U
39638-32-9	bis(2-chloroisopropoxy)Ethane	J
106-44-5	4-Methoxyphenol	J
621-64-7	N-Nitroso-Di-n-Propylamine	J
67-72-1	Hexachloroethane	
98-95-3	Nitrobenzene	
78-59-1	Isophorone	
88-75-5	2-Nitrophenol	
105-67-9	2,4-Dimethoxyphenol	
65-85-0	Benzoic Acid	30,000U
111-91-1	bis(2-Chloroethoxy)Methane	6100U
120-83-2	2,4-Dichlorophenol	
120-82-1	1,2,4-Trichlorobenzene	
91-20-3	Naphthalene	20,000
106-47-8	4-Chloroaniline	6100U
87-68-3	Hexachlorobutadiene	
59-50-7	4-Chloro-3-Methoxyphenol	
91-57-6	2-Methylnaphthalene	10,000
77-47-4	Hexachlorocyclopentadiene	6100U
88-06-2	2,4,6-Trichlorophenol	
95-95-4	2,4,5-Trichlorophenol	30,000U
91-58-7	2-Chloronaphthalene	6100U
88-74-4	2-Nitroaniline	30,000U
131-11-3	Dimethyl Phthalate	6100U
208-96-8	Acenaphthylene	
99-09-2	3-Nitroaniline	30,000U

CAS Number	Compound	ug/l or ug/g (Circle One)
83-32-9	Acenaphthene	6100U
51-28-5	2,4-Dinitrophenol	30,000U
100-02-7	4-Nitrophenol	
132-64-9	Dibenzofuran	6100U
121-14-2	2,4-Dinitrotoluene	
606-20-2	2,5-Dinitrotoluene	
84-66-2	Dienylphthalate	
7005-72-3	4-Chlorophenylphenyl ether	
86-73-7	Fluorene	
100-01-6	4-Nitroaniline	30,000U
534-52-1	4,6-Dinitro-2-Methoxyphenol	
86-30-6	N-Nitrosodiphenylamine (1)	20,000
101-55-3	4-Bromophenylphenyl ether	6100U
118-74-1	Hexachlorobenzene	
87-86-5	Pentachlorophenol	55,000
85-01-8	Phenanthrene	15,000U
120-12-7	Anthracene	6100U
84-74-2	Di-n-Butylphthalate	
206-44-0	Fluoranthene	
129-00-0	Pyrene	
85-68-7	Butylbenzylphthalate	
91-94-1	3,3-Dichlorobenzidine	12,000U
56-55-3	Benzofluoranthene	6100U
117-81-7	bis(2-Ethylhexyl)Phthalate	100,000
218-01-9	Chrysene	6100U
117-84-0	Di-n-Octyl Phthalate	
205-99-2	Benzofluoranthene	J
207-08-9	Benzofluoranthene	J
50-32-8	Benzofluoranthene	
193-39-5	Indeno[1,2,3-cd]Pyrene	
53-70-3	Dibenzofluoranthene	
191-24-2	Benzofluoranthene	

(1)-Cannot be separated from diphenylamine

Laboratory Name ATI  
 Case No 6855 / CAS 2726 E

Sample Number  
1:10  
EJ105

Organics Analysis Data Sheet  
 (Page 3)

RECEIVED MAR 30 1987

Pesticide/PCBs

Concentration: Low Medium (Circle One)  
 Date Extracted/Prepared: 19 FEB 87  
 Date Analyzed: 5 MARCH 87  
 Conc/Dil Factor: 0.1  
 Percent Moisture (decanted) 43.8

GPC Cleanup  Yes  No  
 Separatory Funnel Extraction  Yes  
 Continuous Liquid - Liquid Extraction  Yes

CAS Number		ug/l or ug/Kg (Circle One)
319-84-6	Alpha-BHC	1400M
319-85-7	Beta-BHC	
319-86-8	Delta-BHC	
58-89-9	Gamma-BHC (Lindane)	
76-44-8	Heptachlor	
309-00-2	Aldrin	
1024-57-3	Heptachlor Epoxide	
959-98-8	Endosulfan I	↓
60-57-1	Dieldrin	2800M
72-55-9	4, 4'-DDE	
72-20-8	Endrin	
33213-65-9	Endosulfan II	
72-54-8	4, 4'-DDD	
1031-07-8	Endosulfan Sulfate	↓
50-29-3	4, 4'-DDT	↓
72-43-5	Methoxychlor	1400M
53494-70-5	Endrin Ketone	2800M
57-74-9	Chlordane	1400M
8001-35-2	Toxaphene	2800M
12674-11-2	Aroclor-1016	1400M
11104-28-2	Aroclor-1221	↓
11141-16-5	Aroclor-1232	↓
53469-21-9	Aroclor-1242	↓
12672-29-6	Aroclor-1248	↓
11097-69-1	Aroclor-1254	2800M
11096-82-5	Aroclor-1260	↓

$V_i$  = Volume of extract injected (ul)

$V_s$  = Volume of water extracted (ml)

$W_s$  = Weight of sample extracted (g)

$V_t$  = Volume of total extract (ul)

$V_s$  \_\_\_\_\_ or  $W_s$  30.5  $V_i$  20000  $V_t$  2

Laboratory Name: ATI  
Case No 6855/SAS 2726E

Sample Number  
EJ-105

RECEIVED MAR 30 1987

Organics Analysis Data Sheet  
(Page 4)

Tentatively Identified Compounds

CAS Number	Compound Name	Fraction	RT or Scan Number	Estimated Concentration (ug/l or ug/kg)
1. <u>75-28-5</u>	<u>2-Methyl propanoic acid</u>	<u>VOA</u>	<u>216</u>	<u>88,000 J</u>
2. <u>-</u>	<u>an Acid</u>	<u>↓</u>	<u>613</u>	<u>4,400 J</u>
3.				
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30.				

Laboratory Name: ATI  
 Case No 6855/SAS 2726 E

Sample Number  
EJ105

Organics Analysis Data Sheet  
 (Page 4)

RECEIVED MAR 30 1987

Tentatively Identified Compounds

CAS Number	Compound Name	Fraction	RT or Scan Number	Estimated Concentration (ug/l or ug/kg)
1. —	unknown $\approx$ C <sub>12</sub> hydrocarbon	BNA	1108	60,000 J
2. —	unk. branched hyd. $\approx$ C <sub>10</sub>		1207	78,000 J
3. —	unk. branched phenol		1313	340,000 J
4. —	unk. branched hyd. $\approx$ C <sub>13</sub>		1345	92,000 J
5. —	unk. branched hyd $\approx$ C <sub>17</sub>		1392	320,000 J
6. —	unk. branched phenol		1403	180,000 J
7. —	unk. branched phenol		1411	240,000 J
8. —	unknown		1416	200,000 J
9. —	unk. branched phenol		1442	130,000 J
10. —	unk. phthalate		1449	100,000 J
11. —	unknown		1532	200,000 J
12. —	unknown		1604	280,000 J
13. —	unknown		1614	360,000 J
14. —	unknown		1621	540,000 J
15. —	unk. branched alcohol		1629	220,000 J
16. —	unk. organic acid		1639	440,000 J
17. —	unknown		1653	480,000 J
18. —	unknown	N	1660	140,000 J
19.				
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27 FPD 1561  
 Sample Number  
 EJ-106

RECEIVED MAR 30 1987 Organics Analysis Data Sheet  
 (Page 1)

Laboratory Name: ATI  
 Lab Sample ID No: EJ-106  
 Sample Matrix: SOIL  
 Data Release Authorized By: C. Sites

Case No: 6855/SAS 2726E  
 QC Report No: -  
 Contract No: 6801-7014  
 Date Sample Received: 2/19/87

Volatile Compounds

Concentration: Low Medium (Circle One)  
 Date Extracted/Prepared: 2-24-87  
 Date Analyzed: 2-24-87  
 Conc/Dil Factor: 5g/5g (1.0) pH 7.71  
 Percent Moisture: (Not Decanted) 4.4

CAS Number	Compound	ug/l or ug/Kg (Circle One)
74-87-3	Chloromethane	10U
74-83-9	Bromomethane	
75-01-4	Vinyl Chloride	
75-00-3	Chloroethane	
75-09-2	Methylene Chloride	17B
67-64-1	Acetone	11B
75-15-0	Carbon Disulfide	15U
75-35-4	1, 1-Dichloroethene	
75-34-3	1, 1-Dichloroethane	
156-60-5	Trans-1, 2-Dichloroethene	
67-66-3	Chloroform	
107-06-2	1, 2-Dichloroethane	
78-93-3	2-Butanone	10U
71-55-6	1, 1, 1-Trichloroethane	11J
56-23-5	Carbon Tetrachloride	5U
108-05-4	Vinyl Acetate	10U
75-27-4	Bromodichloromethane	5U

CAS Number	Compound	ug/l or ug/Kg (Circle One)
78-87-5	1, 2-Dichloropropane	5U
10061-02-6	Trans-1, 3-Dichloropropene	
79-01-6	Trichloroethene	
124-48-1	Dibromochloromethane	
79-00-5	1, 1, 2-Trichloroethane	
71-43-2	Benzene	13J
10061-01-5	cis-1, 3-Dichloropropene	5U
110-75-8	2-Chloroethylvinylether	10U
75-25-2	Bromoform	13JB
108-10-1	4-Methyl-2-Pentanone	48U
591-78-6	2-Hexanone	11J
127-18-4	Tetrachloroethene	5U
79-34-5	1, 1, 2, 2-Tetrachloroethane	
108-88-3	Toluene	14JB
108-90-7	Chlorobenzene	5U
100-41-4	Ethylbenzene	13JB
100-42-5	Styrene	5U
	Total Xylenes	112B

Data Reporting Qualifiers

For reporting results to EPA, the following results qualifiers are used. Additional flags or footnotes explaining results are encouraged. However, the definition of each flag must be explicit.

- Value** If the result is a value greater than or equal to the detection limit, report the value
- U** Indicates compound was analyzed for but not detected. Report the minimum detection limit for the sample with the U (e.g., 10U) based on necessary concentration/dilution action. (This is not necessarily the instrument detection limit.) The footnote should read U-Compound was analyzed for but not detected. The number is the minimum attainable detection limit for the sample
- J** Indicates an estimated value. This flag is used either when estimating a concentration for tentatively identified compounds where a 1:1 response is assumed or when the mass spectral data indicated the presence of a compound that meets the identification criteria but the result is less than the specified detection limit but greater than zero (e.g., 10J). If limit of detection is 10 ug/l and a concentration of 3 ug/l is calculated, report as 3J
- C** This flag applies to pesticide parameters where the identification has been confirmed by GC/MS. Single component pesticides  $\geq 10$  ng/l in the final extract should be confirmed by GC/MS
- B** This flag is used when the analyte is found in the blank as well as a sample. It indicates possible/probable blank contamination and warns the data user to take appropriate action
- Other** Other specific flags and footnotes may be required to properly define the results. If used, they must be fully described and such description attached to the data summary report

Laboratory Name ATI  
 Case No: 6855/SAS 2726 E

Sample Number  
EJ 106

Organics Analysis Data Sheet  
 (Page 2)

RECEIVED MAR 30 1987

Semivolatile Compounds

Concentration:  Low Medium (Circle One)  
 Date Extracted/Prepared 2/19/87  
 Date Analyzed: 2/25/87  
 Conc./Dil Factor:  $\frac{1 \text{ mL}}{20 \mu\text{L}} \times 10^3 \times 4 = \times 140$   
 Percent Moisture (Decanted) —

GPC Cleanup  Yes  No  
 Separatory Funnel Extraction  Yes  
 Continuous Liquid - Liquid Extraction  Yes

CAS Number	Compound	ug/l or ug/kg (Circle One)
108-95-2	Phenol	1400 U
111-44-4	bis(2-Chloroethyl)Ether	
95-57-8	2-Chlorophenol	
541-73-1	1,3-Dichlorobenzene	
106-46-7	1,4-Dichlorobenzene	
100-51-6	Benzyl Alcohol	J
95-50-1	1,2-Dichlorobenzene	
95-48-7	2-Methylnaphthalene	
39638-32-9	bis(2-chloroisopropyl)Ether	
106-44-5	4-Methylnaphthalene	
621-64-7	N-Nitroso-Di-n-Propylamine	
67-72-1	Hexachloroethane	
98-95-3	Nitrobenzene	
78-59-1	Isophorone	
88-75-5	2-Nitrophenol	
105-67-9	2,4-Dimethylnaphthalene	71400 U
65-85-0	Benzoic Acid	7000 U, T
111-91-1	bis(2-Chloroethoxy)methane	1400 U
120-83-2	2,4-Dichlorophenol	
120-82-1	1,2,4-Trichlorobenzene	
91-20-3	Naphthalene	
106-47-8	4-Chloroaniline	
87-68-3	Hexachlorocyclopentadiene	
59-50-7	4-Chloro-3-Methylnaphthalene	
91-57-6	2-Methylnaphthalene	
77-47-4	Hexachlorocycloheptadiene	J, T
88-06-2	2,4,6-Trichlorophenol	1400 U
95-95-4	2,4,5-Trichlorophenol	7000 U
91-58-7	2-Chloronaphthalene	1400 U
88-74-4	2-Nitroaniline	7000 U, T
131-11-3	Dimethyl Phthalate	1400 U
208-96-8	Acenaphthylene	J
99-09-2	3-Nitroaniline	7000 U, R

CAS Number	Compound	ug/l or ug/kg (Circle One)
83-32-9	Acenaphthene	1400 U
51-28-5	2,4-Dinitrophenol	7000 U
100-02-7	4-Nitrophenol	J
132-64-9	Dibenzofuran	1400 U
121-14-2	2,4-Dinitrotoluene	
606-20-2	2,6-Dinitrotoluene	
84-66-2	Dibenzylphthalate	
7005-72-3	4-Chlorophenylphenyl ether	J
86-73-7	Fluorene	1400 U
100-01-6	4-Nitroaniline	7000 U, R
534-52-1	4,6-Dinitro-2-Methylnaphthalene	J
86-30-6	N-Nitrosodiphenylamine (1)	1400 U
101-55-3	4-Bromophenylphenyl ether	J
118-74-1	Hexachlorobenzene	J, X
87-86-5	Pentachlorophenol	7000 U
85-01-8	Phenanthrene	1400 U
120-12-7	Anthracene	J
84-74-2	Di-n-Butylphthalate	1100 J
206-44-0	Fluoranthene	240 J
129-00-0	Pyrene	250 J
85-68-7	Butylbenzylphthalate	7000 U
91-94-1	3,3'-Dichlorobenzidine	2800 U, R
56-55-3	Benzofluoranthene	1400 U
117-81-7	bis(2-Ethylhexyl)phthalate	1400 U, T
218-01-9	Chrysene	7000 U
117-84-0	Di-n-Octyl Phthalate	J
205-99-2	Benzobifluoranthene	J
207-08-9	Benzokifluoranthene	
50-32-8	Benzofluorene	
193-39-5	Indeno(1,2,3-cd)pyrene	
53-70-3	Dibenzofluoranthene	J
191-24-2	Benzofluorene	J

(1)-Cannot be separated from diphenylamine

Laboratory Name ATI RECEIVED MAR 30 1987  
 Case No 6855/SAS 2726E

Sample Number  
1.5  
EJ 106

Organics Analysis Data Sheet  
 (Page 3)

Pesticide/PCBs

Concentration: (Low) Medium (Circle One)  
 Date Extracted/Prepared: 19 FEB 87  
 Date Analyzed: 4 MARCH 87  
 Conc/Dil Factor: 0.2  
 Percent Moisture (decanted) 4.44

GPC Cleanup  Yes  No  
 Separatory Funnel Extraction  Yes  
 Continuous Liquid - Liquid Extraction  Yes

CAS Number		ug/l or <u>(ug/Kg)</u> (Circle One)
319-84-6	Alpha-BHC	42m
319-85-7	Beta-BHC	
319-86-8	Delta-BHC	
58-89-9	Gamma-BHC (Lindane)	
76-44-8	Heptachlor	
309-00-2	Aldrin	
1024-57-3	Heptachlor Epoxide	
959-98-8	Endosulfan I	↓
60-57-1	Dieldrin	84m
72-55-9	4,4'-DDE	
72-20-8	Endrin	
33213-65-9	Endosulfan II	
72-54-8	4,4'-DDD	
1031-07-8	Endosulfan Sulfate	
50-29-3	4,4'-ODT	↓
72-43-5	Methoxychlor	420m
53494-70-5	Endrin Ketone	84m
57-74-9	Chlordane	420m
8001-35-2	Toxaphene	840m
12674-11-2	Aroclor-1016	420m
11104-28-2	Aroclor-1221	
11141-16-5	Aroclor-1232	
53469-21-9	Aroclor-1242	
12672-29-6	Aroclor-1248	↓
11097-69-1	Aroclor-1254	840m
11096-82-5	Aroclor-1260	↓

$V_i$  = Volume of extract injected (ul)

$V_s$  = Volume of water extracted (ml)

$W_s$  = Weight of sample extracted (g)

$V_t$  = Volume of total extract (ul)

$V_s$  \_\_\_\_\_ or  $W_s$  30.1  $V_t$  20000  $V_i$  2

Laboratory Name ATI RECEIVED MAR 30 1987  
 Case No 6855 / SAS 2726E

Sample Number  
 1:10  
 EJ 920

Organics Analysis Data Sheet  
 (Page 3)

Pesticide/PCBs

Concentration: Low Medium (Circle One)  
 Date Extracted/Prepared: 19 FEB 87  
 Date Analyzed: 4 MARCH 87  
 Conc/Dil Factor: 0.1  
 Percent Moisture (decanted) 16.2

GPC Cleanup  Yes  No  
 Separatory Funnel Extraction  Yes  
 Continuous Liquid-Liquid Extraction  Yes

CAS Number		ug/l or ug/Kg (Circle One)
319-84-6	Alpha-BHC	94m
319-85-7	Beta-BHC	
319-86-8	Delta-BHC	
58-89-9	Gamma-BHC (Lindane)	
76-44-8	Heptachlor	
309-00-2	Aldrin	
1024-57-3	Heptachlor Epoxide	
959-98-8	Endosulfan I	↓
60-57-1	Dieldrin	190m
72-55-9	4, 4'-DDE	
72-20-8	Endrin	
33213-65-9	Endosulfan II	
72-54-8	4, 4'-DDD	
1031-07-8	Endosulfan Sulfate	
50-29-3	4, 4'-DDT	↓
72-43-5	Methoxychlor	940m
53494-70-5	Endrin Ketone	190m
57-74-9	Chlordane	940m
8001-35-2	Toxaphene	1900m
12674-11-2	Aroclor-1016	940m
11104-28-2	Aroclor-1221	
11141-16-5	Aroclor-1232	
53469-21-9	Aroclor-1242	
12672-29-6	Aroclor-1248	↓
11097-69-1	Aroclor-1254	1900m
11096-82-5	Aroclor-1260	↓

$V_i$  = Volume of extract injected (ul)

$V_s$  = Volume of water extracted (ml)

$W_s$  = Weight of sample extracted (g)

$V_t$  = Volume of total extract (ul)

$V_s$  \_\_\_\_\_ or  $W_s$  30.6  $V_t$  20000  $V_i$  2

Laboratory Name AT7

Case No. 6855/895 2726 E

RECEIVED MAR 30 1987

Sample Number  
EJ 920

Organics Analysis Data Sheet  
(Page 2)

Semivolatile Compounds

Concentration: Low Medium (Circle One)

Date Extracted / Prepared 2/19/87

Date Analyzed: 2/23/87

Conc./Dil Factor: 3.26ml / 32.6g (838) x 10<sup>3</sup> = x 200

Percent Moisture (Decanted) \_\_\_\_\_

GPC Cleanup  Yes  No

Separatory Funnel Extraction  Yes

Continuous Liquid-Liquid Extraction  Yes

CAS Number	Compound	ug/l or ug/Kg (Circle One)
108-95-2	Phenol	2000U
111-44-4	bis(2-Chloroethyl)Ether	
95-57-8	2-Chlorophenol	
541-73-1	1,3-Dichlorobenzene	
106-46-7	1,4-Dichlorobenzene	
100-51-6	Benzyl Alcohol	JFK
95-50-1	1,2-Dichlorobenzene	
95-48-7	2-Methoxyphenol	
39638-32-9	bis(2-chloroisopropyl)Ether	
106-44-5	4-Methoxyphenol	
621-64-7	N-Nitroso-Di-n-Propylamine	
67-72-1	Hexachloroethane	
98-95-3	Nitrobenzene	
78-59-1	Isopropyl	1830J
88-75-5	2-Nitrophenol	2000U
105-67-9	2,4-Dimethoxyphenol	
65-85-0	Benzoic Acid	10,000U
111-91-1	bis(2-Chloroethoxy)Methane	2000U
120-83-2	2,4-Dichlorophenol	
120-82-1	1,2,4-Trichlorobenzene	
91-20-3	Naphthalene	
106-47-8	4-Chloroaniline	RJK
87-68-3	Hexachlorocyclopentadiene	
59-50-7	4-Chloro-3-Methoxyphenol	
91-57-6	2-Methoxynaphthalene	
77-47-4	Hexachlorocyclopentadiene	
88-06-2	2,4,6-Trichlorophenol	
95-95-4	2,4,5-Trichlorophenol	10,000U
91-58-7	2-Chloronaphthalene	2,000U
88-74-4	2-Nitroaniline	10,000U
131-11-3	Dimethyl Phthalate	2,000U
208-96-8	Acenaphthylene	
99-09-2	3-Nitroaniline	10,000U

CAS Number	Compound	ug/l or ug/Kg (Circle One)
83-32-9	Acenaphthene	2000U
51-28-5	2,4-Dinitrophenol	10,000U
100-02-7	4-Nitrophenol	↓
132-64-9	Dibenzofuran	2,000U
121-14-2	2,4-Dinitrotoluene	
606-20-2	2,6-Dinitrotoluene	
84-66-2	Diethylphthalate	
7005-72-3	4-Chlorophenylphenyl ether	
86-73-7	Fluorene	
100-01-6	4-Nitroaniline	10,000U RJK
534-52-1	4,6-Dinitro-2-Methoxyphenol	↓
86-30-6	N-Nitrosodiphenylamine (1)	2,000U RJK
101-55-3	4-Bromophenylphenyl ether	
118-74-1	Hexachlorobenzene	↓
87-86-5	Pentachlorophenol	10,000U
85-01-8	Phenanthrene	1800J
120-12-7	Anthracene	2000U
84-74-2	Di-n-Butylphthalate	1300J
206-44-0	Fluoranthene	1770J
129-00-0	Pyrene	1540J
85-68-7	Butylbenzylphthalate	2000U
91-94-1	3,3-Dichlorobenzidine	4000U
56-55-3	Benz[a]Anthracene	1450J
117-81-7	bis(2-Ethylhexyl)Phthalate	2200
218-01-9	Chrysene	1540J
117-84-0	Di-n-Octyl Phthalate	2000U
205-99-2	Benz[a]Fluoranthene	1550J
207-08-3	Benz[a]Fluoranthene	2000U
50-32-8	Benz[a]Pyrene	1560J
193-39-5	Indeno[1,2,3-cd]Pyrene	2000U
53-70-3	Dibenz[a,h]Anthracene	
191-24-2	Benz[a,h]Perylene	↓

RJK (1) - Cannot be separated from diphenylamine

87FY01562  
 Sample Number  
 EJ 920

RECEIVED MAR 30 1987  
 Organics Analysis Data Sheet  
 (Page 1)

Laboratory Name: ATI  
 Lab Sample ID No: EJ 920  
 Sample Matrix: Soil  
 Data Release Authorized By: C. Sites

Case No: 6855/SAS 2726 E  
 QC Report No: \_\_\_\_\_  
 Contract No: 68017014  
 Date Sample Received: 2/19/87

Volatile Compounds

Concentration: Low Medium (Circle One)  
 Date Extracted/Prepared: 2/24/87  
 Date Analyzed: 2/24/87  
 Conc/Dil Factor: 300 50 (8.33) = 1.2 pH 7.02  
 Percent Moisture: (Not Decanted) 16.2

CAS Number	Compound	ug/l or ug/Kg (Circle One)
74-87-3	Chloromethane	12U
74-83-9	Bromomethane	
75-01-4	Vinyl Chloride	
75-00-3	Chloroethane	
75-09-2	Methylene Chloride	110 B
67-64-1	Acetone	123 B
75-15-0	Carbon Disulfide	6U
75-35-4	1, 1-Dichloroethane	
75-34-3	1, 1-Dichloroethane	
156-60-5	Trans-1, 2-Dichloroethane	
67-66-3	Chloroform	
107-06-2	1, 2-Dichloroethane	12U
78-93-3	2-Butanone	6U
71-55-6	1, 1, 1-Trichloroethane	6U
56-23-5	Carbon Tetrachloride	
108-05-4	Vinyl Acetate	12U
75-27-4	Bromodichloromethane	6U

CAS Number	Compound	ug/l or ug/Kg (Circle One)
78-87-5	1, 2-Dichloropropane	6U
10061-02-6	Trans-1, 3-Dichloropropene	
79-01-6	Trichloroethene	
124-48-1	Dibromochloromethane	
79-00-5	1, 1, 2-Trichloroethane	6U
71-43-2	Benzene	14.71
10061-01-5	cis-1, 3-Dichloropropene	6U
110-75-8	2-Chloroethylvinylether	12U R
75-25-2	Bromoform	6U
108-10-1	4-Methyl-2-Pentanone	12U
591-78-6	2-Hexanone	1.75
127-18-4	Tetrachloroethene	6U
79-34-5	1, 1, 2, 2-Tetrachloroethane	6U
108-88-3	Toluene	15.78 1
108-90-7	Chlorobenzene	6U
100-41-4	Ethylbenzene	13.78 1
100-42-5	Styrene	6U
	Total Xylenes	71.8 1

Data Reporting Qualifiers

For reporting results to EPA, the following results qualifiers are used. Additional flags or footnotes explaining results are encouraged. However, the definition of each flag must be explicit.

- U** If the result is a value greater than or equal to the detection limit, report the value
- U** Indicates compound was analyzed for but not detected. Report the minimum detection limit for the sample with the U (e.g., 10U) based on necessary concentration/dilution action. (This is not necessary if the instrument detection limit.) The footnote should read: U-Compound was analyzed for but not detected. The number is the minimum attainable detection limit for the sample
- U** Indicates an estimated value. This flag is used either when estimating a concentration for tentatively identified compounds where a 1:1 response is assumed or when the mass spectral data indicated the presence of a compound that meets the identification criteria but the result is less than the specified detection limit but greater than zero (e.g., 10U). If limit of detection is 10 ug/l and a concentration of 3 ug/l is calculated, report as 3U

- C** This flag applies to pesticide parameters where the identification has been confirmed by GC/MS. Single component pesticides  $\geq 10$  ng/l in the final extract should be confirmed by GC/MS
- B** This flag is used when the analyte is found in the blank as well as a sample. It indicates possible/probable blank contamination and warns the data user to take appropriate action
- Other** Other specific flags and footnotes may be required to properly define the results. If used, they must be fully described and such description attached to the data summary report

Laboratory Name: ATI RECEIVED MAR 30 1987  
 Case No 6855 / SAS 2726 E

Sample Number  
EJ 106

Organics Analysis Data Sheet  
 (Page 4)

Tentatively Identified Compounds

CAS Number	Compound Name	Fraction	RT or Scan Number	Estimated Concentration (ug/l or ug/kg)
1. <u>—</u>	<u>Unknown</u>	<u>BNA</u>	<u>623</u>	<u>1100J</u>
2.				
3.				
4.				
5.				
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Laboratory Name: ATI  
 Case No 6855 / SAS 2726 E RECEIVED MAR 30 1987

Sample Number  
EJ-106

Organics Analysis Data Sheet  
 (Page 4)

Tentatively Identified Compounds

CAS Number	Compound Name	Fraction	RT or Scan Number	Estimated Concentration (ug/l or ug/kg)
1. —	NONE FOUND	VOA	—	—
2.				
3.				
4.				
5.				
6.				
7.				
8.				
9.				
10.				
11.				
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28.				
29.				
30.				

Laboratory Name: ATT  
 Case No 6855/87AS 2726 E

RECEIVED MAR 30 1987

Sample Number  
EJ920

Organics Analysis Data Sheet  
 (Page 4)

Tentatively Identified Compounds

CAS Number	Compound Name	Fraction	RT or Scan Number	Estimated Concentration (ug/l or ug/kg)
1. —	Unknown Xylene	BNA	418	1000 J
2. —	Unknown Xylene	"	459	800 J
3. —	Unknown branched Phenol		1233	1200 J
4. —	Unknown		1548	2600 J
5. —	Unknown		1660	1400 J
6. —	Unknown		1768	13000 J
7. —	Unknown		2026	400 J
8. —	Unknown		2199	3200 J
9. —	Unknown		2268	3000 J
10. —	Unknown	↓	2387	2600 J
11.				
12.				
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14.				
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# ecology and environment, inc.

111 WEST JACKSON BLVD., CHICAGO, ILLINOIS 60604, TEL. 312-663-9415

International Specialists in the Environment

Date Received for Review: 3/30/87 Date Review Completed: 3/30/87

TO: Julie Kaiser

FROM: Zena Gold-Kaufman

SUBJECT: MOBIL CHEM  
FINOTTO  
FOS-8612-012

Sample Description: Case # 6855

4 soil samples-organics

Project Data Status: complete\*

## FIT Date Review Findings:

several organic hits - see  
Tom Clynt's review

## Additional Comments:

none

Book No. 6

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UNITED STATES ENVIRONMENTAL PROTECTION AGENCY  
REGION V

DATE: 3/26/87 RECEIVED MAR 30 1987  
TO: Data User: \_\_\_\_\_ FIT \_\_\_\_\_

FROM: Curtis Ross, Director (5SCRL) *Patrick J. Chumilla*  
Central Regional Laboratory

SUBJECT: Review of Region V CLP Data  
Received for Review on 3/6/87

We have reviewed the data for the following case(s).

SITE NAME Mobil Chem SMO Case No. 6855/2726 SAS

EPA DATA SET No. SF3778 No. OF SAMPLES 4 D.U./ACTIVITY NUMBERS Y905/C72100

CRL No.: 87FP01559-87FP0/S62

SMO Traffic No. EJ104-EJ106; EJ920

CLP Laboratory: ATI Hrs. Required for Review: 12

Following are our findings.

- A. Time Acceptable
- B. Surrogates Acceptable
- C. MS/MSD EJ105 VOA was run at medium level and has 4/10 outliers
- D. Lab Blanks
  - Pest Acceptable
  - 2/19/87 SV Dibutyl phthalate 780mg/kg and 5 TICs
  - 2/24/87 VOA Methylene chloride 6 mg/L 5 other HSL < CRDL
  - 2/25/87 VOA Methylene chloride 1000 mg/L
  - Chloroform 250
  - Ethylbenzene 300
  - Total xylenes 380
  - 3 other HSL < CRDL
- E. Tuning Acceptable
- F. Calibration Acceptable, see calibration outlier forms
- G. Pesticides Acceptable

- ( ) Data are acceptable for use.
- ( X ) Data are acceptable for use with qualifications noted above.
- ( ) Data are preliminary - pending verification by contractor lab.
- ( ) Data are unacceptable.

cc: Dr. Alfred Haerber/Joan Fisk/Gary Ward. EPA Support Services.  
Ross K. Robeson, EMSL - LasVegas  
Don Trees, CLP/Sample Management Office

*Reviewed  
25 Mar 87  
Tom Lynn*

USER'S INFORMATION FORM

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The follow samples have HRS hits that can be used. Some hits were found in the blanks but because they are 10 x the value found in the lab blank, they can be used.

EJ 104	VOA	benzene	41 mg/L , this is a background sample.
EJ 105	VOA	toluene	
		ethyl benzene, total xylenes	
	SV	1,2 dichlorobenzene	
		<del>naphthalene</del> <i>TC NAPHTHALENE</i>	
		2-methyl-naphthalene	
		3-nitroaniline	
		n-nitrosodiphenylamine	
		pentachlorophenol	
		bis-(2-ethyl hexyl)phthalate	
EJ 106	SV	bis-(2-ethyl hexyl)phthalate	
ET 920		No usable hits	

Team leader, please check EJ 104 and evaluate if this site was a true background. This compound could be in the soil or it could be a lab artifact.

